

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041625\
 Data File : BM049957.D
 Acq On : 16 Apr 2025 18:03
 Operator : RC/JU
 Sample : Q1808-01MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OILY-SOIL-PILEMSD

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 04/17/2025
 Supervised By :Jagrut Upadhyay 04/17/2025

Quant Time: Apr 16 18:40:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	369676	20.000	ng	-0.01
21) Naphthalene-d8	10.563	136	1237619	20.000	ng	-0.01
39) Acenaphthene-d10	14.410	164	766997	20.000	ng	-0.01
64) Phenanthrene-d10	17.157	188	1513554	20.000	ng	0.00
76) Chrysene-d12	21.403	240	1581516	20.000	ng	0.00
86) Perylene-d12	24.403	264	1732464	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	1531602	70.353	ng	0.00
7) Phenol-d6	6.951	99	1963013	72.473	ng	0.00
23) Nitrobenzene-d5	8.928	82	1120213	50.403	ng	0.00
42) 2,4,6-Tribromophenol	15.904	330	804142	72.080	ng	0.00
45) 2-Fluorobiphenyl	13.033	172	2699268	47.722	ng	-0.01
79) Terphenyl-d14	19.786	244	3881366	45.993	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.252	88	331801	37.008	ng	99
3) Pyridine	3.652	79	886022	37.613	ng	99
4) n-Nitrosodimethylamine	3.558	42	357717	39.903	ng	98
6) Aniline	7.098	93	693846	29.895	ng	100
8) 2-Chlorophenol	7.340	128	909259	40.755	ng	100
9) Benzaldehyde	6.910	77	450756	32.429	ng	99
10) Phenol	6.975	94	1102282	41.702	ng	100
11) bis(2-Chloroethyl)ether	7.193	93	848288	40.925	ng	99
12) 1,3-Dichlorobenzene	7.657	146	1082747	41.049	ng	99
13) 1,4-Dichlorobenzene	7.804	146	1089518	41.051	ng	99
14) 1,2-Dichlorobenzene	8.122	146	1044648	41.788	ng	99
15) Benzyl Alcohol	8.010	79	677425	39.717	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.293	45	972620	42.099	ng	99
17) 2-Methylphenol	8.222	107	689259	42.315	ng	100
18) Hexachloroethane	8.845	117	371699	40.254	ng	96
19) n-Nitroso-di-n-propyla...	8.575	70	598167	39.873	ng	100
20) 3+4-Methylphenols	8.551	107	908212	40.897	ng	98
22) Acetophenone	8.592	105	1208718	40.186	ng	# 99
24) Nitrobenzene	8.969	77	866159	40.473	ng	96
25) Isophorone	9.498	82	1591002	42.732	ng	99
26) 2-Nitrophenol	9.681	139	475156	41.890	ng	98
27) 2,4-Dimethylphenol	9.751	122	763996	59.433	ng	98
28) bis(2-Chloroethoxy)met...	9.975	93	1007527	40.421	ng	99
29) 2,4-Dichlorophenol	10.228	162	862412	40.360	ng	99
30) 1,2,4-Trichlorobenzene	10.428	180	1029854	41.662	ng	99
31) Naphthalene	10.616	128	2515800	39.560	ng	99
32) Benzoic acid	9.892	122	458701m	32.206	ng	
33) 4-Chloroaniline	10.734	127	454385	20.235	ng	98
34) Hexachlorobutadiene	10.898	225	656410	42.929	ng	100
35) Caprolactam	11.510	113	230307	37.580	ng	99
36) 4-Chloro-3-methylphenol	11.869	107	732863	39.155	ng	98
37) 2-Methylnaphthalene	12.228	142	1627295	36.319	ng	99
38) 1-Methylnaphthalene	12.451	142	1707995	39.128	ng	100
40) 1,2,4,5-Tetrachloroben...	12.604	216	1139380	42.462	ng	98
41) Hexachlorocyclopentadiene	12.580	237	1055285	112.235	ng	100
43) 2,4,6-Trichlorophenol	12.845	196	726527	42.550	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041625\
 Data File : BM049957.D
 Acq On : 16 Apr 2025 18:03
 Operator : RC/JU
 Sample : Q1808-01MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OILY-SOIL-PILEMSD

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 04/17/2025
 Supervised By :Jagrut Upadhyay 04/17/2025

Quant Time: Apr 16 18:40:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.933	196	784458	42.277	ng	98
46) 1,1'-Biphenyl	13.245	154	2348328	41.187	ng	100
47) 2-Chloronaphthalene	13.286	162	1858097	41.463	ng	99
48) 2-Nitroaniline	13.492	65	440557	42.494	ng	96
49) Acenaphthylene	14.133	152	3001811	45.985	ng	100
50) Dimethylphthalate	13.869	163	2149959	39.733	ng	99
51) 2,6-Dinitrotoluene	13.992	165	471424	41.583	ng	94
52) Acenaphthene	14.474	154	1690361	40.710	ng	99
53) 3-Nitroaniline	14.322	138	379621	34.646	ng	96
54) 2,4-Dinitrophenol	14.533	184	458345	58.011	ng	97
55) Dibenzofuran	14.816	168	2747091	39.576	ng	98
56) 4-Nitrophenol	14.657	139	807322	82.681	ng	100
57) 2,4-Dinitrotoluene	14.780	165	656603	43.414	ng	97
58) Fluorene	15.463	166	2280971	41.340	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.045	232	657445	40.425	ng	97
60) Diethylphthalate	15.233	149	1977276	38.104	ng	99
61) 4-Chlorophenyl-phenyle...	15.457	204	1209434	40.885	ng	99
62) 4-Nitroaniline	15.486	138	457581	40.382	ng	99
63) Azobenzene	15.751	77	1868965	42.069	ng	97
65) 4,6-Dinitro-2-methylph...	15.551	198	309039	32.402	ng	97
66) n-Nitrosodiphenylamine	15.668	169	1863778	41.148	ng	100
67) 4-Bromophenyl-phenylether	16.351	248	735992	41.888	ng	96
68) Hexachlorobenzene	16.468	284	859631	42.653	ng	98
69) Atrazine	16.627	200	658237	56.048	ng	98
70) Pentachlorophenol	16.815	266	1175579	79.229	ng	99
71) Phenanthrene	17.204	178	4014857	48.384	ng	100
72) Anthracene	17.292	178	3713619	45.413	ng	100
73) Carbazole	17.562	167	3182056	41.314	ng	100
74) Di-n-butylphthalate	18.127	149	3483798	39.270	ng	99
75) Fluoranthene	19.221	202	5911614	57.942	ng	99
77) Benzidine	19.409	184	2155270	102.725	ng	100
78) Pyrene	19.580	202	5867613	53.834	ng	99
80) Butylbenzylphthalate	20.480	149	1606238	39.219	ng	96
81) Benzo(a)anthracene	21.380	228	5295288	50.380	ng	99
82) 3,3'-Dichlorobenzidine	21.303	252	1323105	33.334	ng	99
83) Chrysene	21.445	228	4847226	48.508	ng	99
84) Bis(2-ethylhexyl)phtha...	21.303	149	2399707	40.216	ng	97
85) Di-n-octyl phthalate	22.433	149	4184046	40.081	ng	# 100
87) Indeno(1,2,3-cd)pyrene	27.803	276	5924472	45.876	ng	98
88) Benzo(b)fluoranthene	23.462	252	5347993	48.334	ng	99
89) Benzo(k)fluoranthene	23.521	252	4960054	46.259	ng	99
90) Benzo(a)pyrene	24.268	252	5005538	51.482	ng	99
91) Dibenzo(a,h)anthracene	27.850	278	4545734	42.735	ng	99
92) Benzo(g,h,i)perylene	28.868	276	4698084	43.222	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

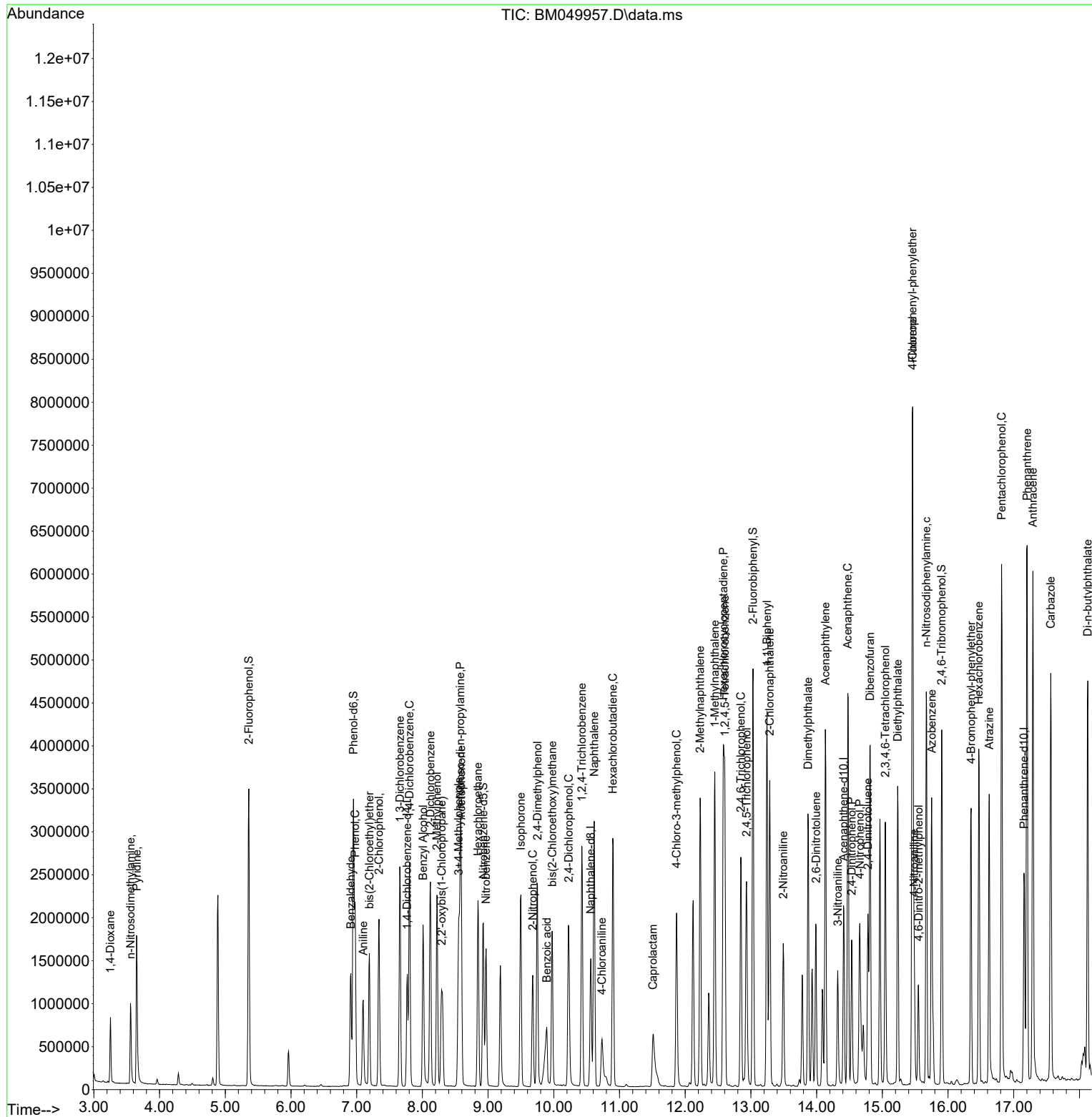
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041625\
Data File : BM049957.D
Acq On : 16 Apr 2025 18:03
Operator : RC/JU
Sample : Q1808-01MSD
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
OILY-SOIL-PILEMSD

Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 04/17/2025
Supervised By :Jagrut Upadhyay 04/17/2025

Quant Time: Apr 16 18:40:39 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 09 04:00:55 2025
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041625\
Data File : BM049957.D
Acq On : 16 Apr 2025 18:03
Operator : RC/JU
Sample : Q1808-01MSD
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
OILY-SOIL-PILEMSD

Quant Time: Apr 16 18:40:39 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 09 04:00:55 2025
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 04/17/2025
Supervised By :Jagrut Upadhyay 04/17/2025

