

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
 Data File : BM047251.D
 Acq On : 19 Aug 2024 13:42
 Operator : RC/JU
 Sample : PB162757BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162757BSD

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/20/2024
 Supervised By :mohammad ahmed 08/21/2024

Quant Time: Aug 19 14:49:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	188275	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	791375	20.000	ng	-0.01
39) Acenaphthene-d10	14.027	164	557332	20.000	ng	0.00
64) Phenanthrene-d10	16.792	188	1252425	20.000	ng	0.00
76) Chrysene-d12	21.033	240	1245983	20.000	ng	0.00
86) Perylene-d12	23.738	264	1572179	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1424206	135.166	ng	0.00
7) Phenol-d6	6.581	99	1895913	130.584	ng	0.00
23) Nitrobenzene-d5	8.516	82	1231431	78.353	ng	0.00
42) 2,4,6-Tribromophenol	15.539	330	920269	142.405	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	2950276	81.655	ng	-0.01
79) Terphenyl-d14	19.462	244	5506473	94.697	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.022	88	113428	25.981	ng	96
3) Pyridine	3.399	79	370799	28.964	ng	97
4) n-Nitrosodimethylamine	3.316	42	205628	36.764	ng	99
6) Aniline	6.710	93	383691	27.960	ng	99
8) 2-Chlorophenol	6.945	128	598876	52.112	ng	97
9) Benzaldehyde	6.528	77	163403	17.223	ng	99
10) Phenol	6.610	94	702697	48.505	ng	98
11) bis(2-Chloroethyl)ether	6.816	93	529915	42.713	ng	95
12) 1,3-Dichlorobenzene	7.251	146	593110	43.431	ng	97
13) 1,4-Dichlorobenzene	7.398	146	605106	43.763	ng	98
14) 1,2-Dichlorobenzene	7.710	146	595731	45.664	ng	99
15) Benzyl Alcohol	7.616	79	535220	51.797	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.887	45	672110	42.218	ng	99
17) 2-Methylphenol	7.828	107	484486	50.128	ng	99
18) Hexachloroethane	8.416	117	222857	41.610	ng	94
19) n-Nitroso-di-n-propyla...	8.175	70	421681	43.118	ng	98
20) 3+4-Methylphenols	8.157	107	670847	49.649	ng	98
22) Acetophenone	8.187	105	752379	39.386	ng	98
24) Nitrobenzene	8.557	77	652013	41.200	ng	96
25) Isophorone	9.081	82	1219558	44.567	ng	98
26) 2-Nitrophenol	9.257	139	342308	48.841	ng	96
27) 2,4-Dimethylphenol	9.339	122	441628	53.793	ng	100
28) bis(2-Chloroethoxy)met...	9.563	93	748328	43.460	ng	100
29) 2,4-Dichlorophenol	9.792	162	640100	52.547	ng	99
30) 1,2,4-Trichlorobenzene	9.992	180	623817	45.252	ng	99
31) Naphthalene	10.169	128	1725748	43.775	ng	99
32) Benzoic acid	9.522	122	486844	51.041	ng	96
33) 4-Chloroaniline	10.298	127	412674	27.220	ng	98
34) Hexachlorobutadiene	10.457	225	363591	45.061	ng	100
35) Caprolactam	11.104	113	203814m	48.179	ng	
36) 4-Chloro-3-methylphenol	11.451	107	637315	49.126	ng	99
37) 2-Methylnaphthalene	11.798	142	1272582	45.337	ng	98
38) 1-Methylnaphthalene	12.022	142	1186326	42.764	ng	98
40) 1,2,4,5-Tetrachloroben...	12.186	216	663901	43.673	ng	98
41) Hexachlorocyclopentadiene	12.157	237	473446	89.192	ng	98
43) 2,4,6-Trichlorophenol	12.445	196	541460	50.113	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
 Data File : BM047251.D
 Acq On : 19 Aug 2024 13:42
 Operator : RC/JU
 Sample : PB162757BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162757BSD

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/20/2024
 Supervised By :mohammad ahmed 08/21/2024

Quant Time: Aug 19 14:49:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.527	196	567752	47.326	ng	96
46) 1,1'-Biphenyl	12.851	154	1586071	39.974	ng	100
47) 2-Chloronaphthalene	12.886	162	1342415	43.287	ng	98
48) 2-Nitroaniline	13.110	65	443955	45.696	ng	98
49) Acenaphthylene	13.745	152	2204793	48.232	ng	100
50) Dimethylphthalate	13.510	163	1805626	46.388	ng	100
51) 2,6-Dinitrotoluene	13.627	165	416728	45.799	ng	87
52) Acenaphthene	14.092	154	1275241	43.979	ng	98
53) 3-Nitroaniline	13.957	138	357076	39.163	ng	96
54) 2,4-Dinitrophenol	14.180	184	550923	101.021	ng	92
55) Dibenzofuran	14.433	168	2087079	45.567	ng	98
56) 4-Nitrophenol	14.304	139	774486	98.507	ng	99
57) 2,4-Dinitrotoluene	14.427	165	573934	47.256	ng	97
58) Fluorene	15.092	166	1718229	45.466	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.674	232	529330	53.688	ng	93
60) Diethylphthalate	14.892	149	1818603	45.752	ng	98
61) 4-Chlorophenyl-phenyle...	15.092	204	834673	46.392	ng	96
62) 4-Nitroaniline	15.133	138	430471	47.923	ng	94
63) Azobenzene	15.392	77	1610976	42.744	ng	94
65) 4,6-Dinitro-2-methylph...	15.204	198	372358	50.780	ng	91
66) n-Nitrosodiphenylamine	15.315	169	1539154	45.119	ng	100
67) 4-Bromophenyl-phenylether	15.992	248	562507	46.609	ng	96
68) Hexachlorobenzene	16.098	284	651895	45.101	ng	98
69) Atrazine	16.292	200	516465	50.312	ng	98
70) Pentachlorophenol	16.457	266	930264	91.815	ng	99
71) Phenanthrene	16.833	178	2889868	45.977	ng	100
72) Anthracene	16.927	178	2909916	47.580	ng	100
73) Carbazole	17.209	167	2910966	46.617	ng	99
74) Di-n-butylphthalate	17.792	149	3370294	46.385	ng	99
75) Fluoranthene	18.868	202	3662035	47.900	ng	99
77) Benzidine	19.074	184	972955	46.046	ng	99
78) Pyrene	19.239	202	3872962	43.533	ng	99
80) Butylbenzylphthalate	20.174	149	1685588	46.426	ng	95
81) Benzo(a)anthracene	21.021	228	3904446	47.203	ng	100
82) 3,3'-Dichlorobenzidine	20.956	252	1267661	48.982	ng	99
83) Chrysene	21.080	228	3667613	46.734	ng	100
84) Bis(2-ethylhexyl)phtha...	20.974	149	2306020	44.786	ng #	99
85) Di-n-octyl phthalate	22.009	149	4339330	49.582	ng	98
87) Indeno(1,2,3-cd)pyrene	26.756	276	5410859	53.245	ng	98
88) Benzo(b)fluoranthene	22.897	252	4408520	47.218	ng	99
89) Benzo(k)fluoranthene	22.956	252	4184991	47.455	ng	99
90) Benzo(a)pyrene	23.621	252	4186037	51.757	ng	98
91) Dibenzo(a,h)anthracene	26.809	278	4361715	53.036	ng	99
92) Benzo(g,h,i)perylene	27.691	276	4213855	49.351	ng	99

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

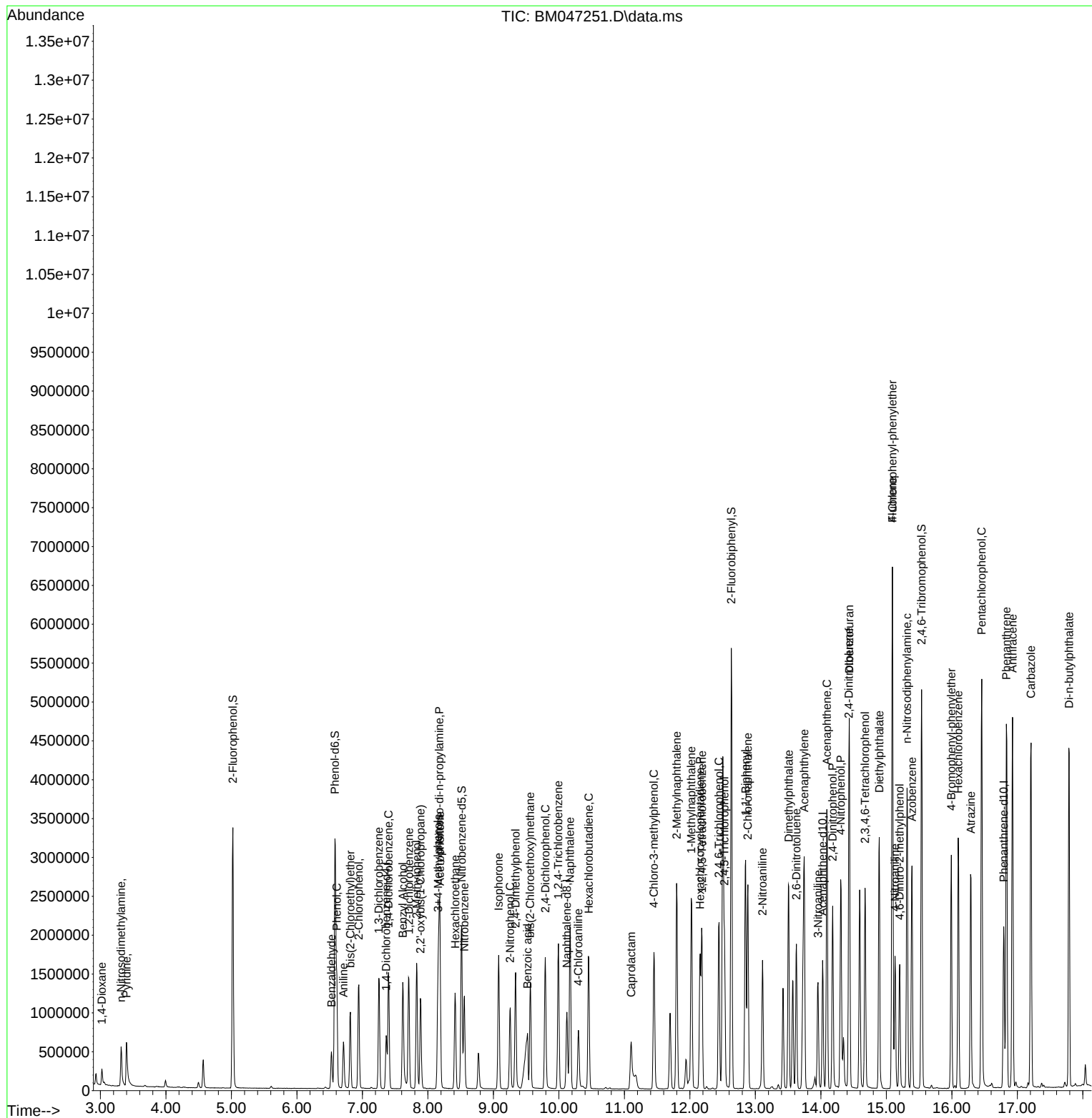
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
Data File : BM047251.D
Acq On : 19 Aug 2024 13:42
Operator : RC/JU
Sample : PB162757BSD
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB162757BSD

Quant Time: Aug 19 14:49:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Aug 11 23:47:58 2024
Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 08/20/2024
Supervised By :mohammad ahmed 08/21/2024



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
Data File : BM047251.D
Acq On : 19 Aug 2024 13:42
Operator : RC/JU
Sample : PB162757BSD
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB162757BSD

Manual Integrations
APPROVED

Reviewed By :Jagrut Upadhyay 08/20/2024
Supervised By :mohammad ahmed 08/21/2024

Quant Time: Aug 19 14:49:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Aug 11 23:47:58 2024
Response via : Initial Calibration

