

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090524\
 Data File : BF139259.D
 Acq On : 05 Sep 2024 15:49
 Operator : RC/JU
 Sample : P2403-02
 Misc : LOQ-SOIL-5 ng
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-SOIL-02-QT2-2024

Quant Time: Sep 05 16:46:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 16:11:43 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	107568	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	390234	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	215486	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	392752	20.000	ng	0.00
76) Chrysene-d12	14.045	240	214033	20.000	ng	0.00
86) Perylene-d12	15.521	264	183805	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	755956	117.564	ng	0.00
7) Phenol-d6	6.510	99	995472	115.966	ng	0.00
23) Nitrobenzene-d5	7.445	82	643189	85.745	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	220980	118.758	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1174236	84.117	ng	0.00
79) Terphenyl-d14	12.998	244	1097832	82.816	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.675	88	18810	6.282	ng	99
3) Pyridine	3.458	79	42053	5.818	ng	96
4) n-Nitrosodimethylamine	3.369	42	27474	5.853	ng	99
6) Aniline	6.546	93	59881	7.586	ng	97
8) 2-Chlorophenol	6.669	128	43068	6.463	ng	87
9) Benzaldehyde	6.434	77	38278	7.367	ng	98
10) Phenol	6.522	94	58273	6.436	ng	# 70
11) bis(2-Chloroethyl)ether	6.622	93	41896	6.459	ng	100
12) 1,3-Dichlorobenzene	6.828	146	46838	6.210	ng	99
13) 1,4-Dichlorobenzene	6.904	146	46940	6.217	ng	99
14) 1,2-Dichlorobenzene	7.057	146	44500	6.320	ng	99
15) Benzyl Alcohol	7.022	79	47571	7.599	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.157	45	81846	6.628	ng	99
17) 2-Methylphenol	7.128	107	34029	6.251	ng	97
18) Hexachloroethane	7.398	117	16310	6.049	ng	98
19) n-Nitroso-di-n-propyla...	7.287	70	32001	6.473	ng	98
20) 3+4-Methylphenols	7.281	107	45081	6.571	ng	97
22) Acetophenone	7.287	105	61143	6.577	ng	99
24) Nitrobenzene	7.463	77	49835	6.259	ng	99
25) Isophorone	7.698	82	82118	6.403	ng	99
26) 2-Nitrophenol	7.781	139	18517	5.922	ng	97
27) 2,4-Dimethylphenol	7.816	122	29023	6.508	ng	98
28) bis(2-Chloroethoxy)met...	7.916	93	48935	6.404	ng	99
29) 2,4-Dichlorophenol	8.016	162	33545	6.282	ng	98
30) 1,2,4-Trichlorobenzene	8.110	180	38693	6.345	ng	99
31) Naphthalene	8.187	128	126879	6.420	ng	99
32) Benzoic acid	7.863	122	13733	3.381	ng	98
33) 4-Chloroaniline	8.234	127	47824	6.989	ng	98
34) Hexachlorobutadiene	8.310	225	23300	6.029	ng	98
35) Caprolactam	8.563	113	9169	5.547	ng	90
36) 4-Chloro-3-methylphenol	8.710	107	35388	6.082	ng	95
37) 2-Methylnaphthalene	8.881	142	80749	6.531	ng	98
38) 1-Methylnaphthalene	8.981	142	74968	6.165	ng	100
40) 1,2,4,5-Tetrachloroben...	9.045	216	38856	6.373	ng	98
41) Hexachlorocyclopentadiene	9.034	237	13099	6.685	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	23951	6.004	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	24688	5.906	ng	98
46) 1,1'-Biphenyl	9.345	154	107383	6.726	ng	100
47) 2-Chloronaphthalene	9.369	162	76448	6.341	ng	97
48) 2-Nitroaniline	9.463	65	24848	5.984	ng	99
49) Acenaphthylene	9.781	152	125893	6.934	ng	99
50) Dimethylphthalate	9.639	163	85950	6.581	ng	99
51) 2,6-Dinitrotoluene	9.698	165	17357	5.858	ng	96
52) Acenaphthene	9.957	154	72970	6.059	ng	95
53) 3-Nitroaniline	9.869	138	19014	6.023	ng	98
54) 2,4-Dinitrophenol	9.975	184	2476	8.971	ng	# 45
55) Dibenzofuran	10.128	168	115051	6.663	ng	99
56) 4-Nitrophenol	10.022	139	11283	4.842	ng	93
57) 2,4-Dinitrotoluene	10.098	165	21852	5.794	ng	96
58) Fluorene	10.469	166	90059	6.728	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	20291	6.167	ng	97
60) Diethylphthalate	10.339	149	82751	6.438	ng	98
61) 4-Chlorophenyl-phenyle...	10.463	204	43258	6.701	ng	98
62) 4-Nitroaniline	10.475	138	17275	5.749	ng	97
63) Azobenzene	10.622	77	86161	6.360	ng	99
65) 4,6-Dinitro-2-methylph...	10.504	198	5977	3.294	ng	98
66) n-Nitrosodiphenylamine	10.581	169	76017	6.649	ng	97
67) 4-Bromophenyl-phenylether	10.957	248	24368	6.395	ng	95
68) Hexachlorobenzene	11.016	284	28094	6.440	ng	99
69) Atrazine	11.104	200	20724	7.857	ng	96
70) Pentachlorophenol	11.210	266	10234	4.061	ng	96
71) Phenanthrene	11.433	178	124829	6.804	ng	99
72) Anthracene	11.486	178	124646	6.954	ng	98
73) Carbazole	11.639	167	108008	6.448	ng	99
74) Di-n-butylphthalate	11.975	149	107247	6.253	ng	98
75) Fluoranthene	12.622	202	122911	6.682	ng	99
77) Benzidine	12.745	184	34855	8.088	ng	97
78) Pyrene	12.851	202	131028	6.337	ng	98
80) Butylbenzylphthalate	13.469	149	28404	5.836	ng	99
81) Benzo(a)anthracene	14.033	228	90356	6.611	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	20971	5.778	ng	99
83) Chrysene	14.074	228	77891	6.106	ng	100
84) Bis(2-ethylhexyl)phtha...	14.033	149	28992	5.075	ng	99
85) Di-n-octyl phthalate	14.645	149	37223	3.403	ng	# 99
87) Indeno(1,2,3-cd)pyrene	16.992	276	54337	4.732	ng	97
88) Benzo(b)fluoranthene	15.086	252	57150	5.399	ng	98
89) Benzo(k)fluoranthene	15.116	252	68354	7.202	ng	98
90) Benzo(a)pyrene	15.457	252	54039	6.012	ng	100
91) Dibenzo(a,h)anthracene	17.015	278	42866	4.530	ng	97
92) Benzo(g,h,i)perylene	17.439	276	42604	4.362	ng	97

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

