

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090524\
 Data File : BF139262.D
 Acq On : 05 Sep 2024 17:25
 Operator : RC/JU
 Sample : P2403-01
 Misc : LOD-MDL-SOIL-8 ng
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT2-2024

Quant Time: Sep 05 17:54:18 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	113163	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	417052	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	227568	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	410152	20.000	ng	0.00
76) Chrysene-d12	14.045	240	205994	20.000	ng	0.00
86) Perylene-d12	15.521	264	186028	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	875864	129.477	ng	0.00
7) Phenol-d6	6.516	99	1166957	129.222	ng	0.00
23) Nitrobenzene-d5	7.451	82	755933	94.295	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	257187	130.878	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1322279	89.693	ng	0.00
79) Terphenyl-d14	12.998	244	1109321	86.949	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.675	88	27685	8.788	ng	96
3) Pyridine	3.457	79	63733	8.381	ng	98
4) n-Nitrosodimethylamine	3.369	42	42988	8.706	ng	97
6) Aniline	6.551	93	90958	10.953	ng	97
8) 2-Chlorophenol	6.669	128	65300	9.314	ng	86
9) Benzaldehyde	6.440	77	58628	10.725	ng	99
10) Phenol	6.528	94	90788	9.531	ng	90
11) bis(2-Chloroethyl)ether	6.622	93	65477	9.595	ng	97
12) 1,3-Dichlorobenzene	6.828	146	72283	9.109	ng	99
13) 1,4-Dichlorobenzene	6.904	146	74014	9.319	ng	97
14) 1,2-Dichlorobenzene	7.057	146	67518	9.115	ng	100
15) Benzyl Alcohol	7.022	79	69293	10.522	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.157	45	122861	9.457	ng	99
17) 2-Methylphenol	7.128	107	51548	9.001	ng	98
18) Hexachloroethane	7.398	117	24698	8.707	ng	98
19) n-Nitroso-di-n-propyla...	7.292	70	48805	9.384	ng	99
20) 3+4-Methylphenols	7.281	107	68090	9.434	ng	92
22) Acetophenone	7.292	105	93337	9.394	ng	94
24) Nitrobenzene	7.469	77	74583	8.765	ng	99
25) Isophorone	7.698	82	128681	9.389	ng	99
26) 2-Nitrophenol	7.781	139	29221	8.744	ng	99
27) 2,4-Dimethylphenol	7.816	122	37768	7.924	ng	99
28) bis(2-Chloroethoxy)met...	7.916	93	76029	9.311	ng	97
29) 2,4-Dichlorophenol	8.022	162	52216	9.150	ng	98
30) 1,2,4-Trichlorobenzene	8.110	180	59084	9.065	ng	98
31) Naphthalene	8.192	128	192152	9.098	ng	99
32) Benzoic acid	7.869	122	20023	4.612	ng	97
33) 4-Chloroaniline	8.234	127	72546	9.920	ng	98
34) Hexachlorobutadiene	8.310	225	35543	8.606	ng	99
35) Caprolactam	8.569	113	14420	8.162	ng	98
36) 4-Chloro-3-methylphenol	8.710	107	54658	8.789	ng	95
37) 2-Methylnaphthalene	8.881	142	122180	9.246	ng	99
38) 1-Methylnaphthalene	8.981	142	112995	8.695	ng	97
40) 1,2,4,5-Tetrachloroben...	9.045	216	59823	9.291	ng	98
41) Hexachlorocyclopentadiene	9.034	237	17426	8.422	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	37119	8.811	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	36417	8.249	ng	97
46) 1,1'-Biphenyl	9.345	154	159559	9.463	ng	100
47) 2-Chloronaphthalene	9.369	162	113432	8.910	ng	98
48) 2-Nitroaniline	9.463	65	37665	8.589	ng	99
49) Acenaphthylene	9.786	152	188521	9.832	ng	100
50) Dimethylphthalate	9.639	163	125745	9.117	ng	100
51) 2,6-Dinitrotoluene	9.698	165	26758	8.551	ng	98
52) Acenaphthene	9.957	154	111484	8.766	ng	98
53) 3-Nitroaniline	9.869	138	28677	8.602	ng	99
54) 2,4-Dinitrophenol	9.975	184	4881	10.223	ng	# 69
55) Dibenzofuran	10.128	168	172032	9.435	ng	98
56) 4-Nitrophenol	10.016	139	17131	6.962	ng	98
57) 2,4-Dinitrotoluene	10.098	165	33719	8.466	ng	99
58) Fluorene	10.469	166	132253	9.355	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	30327	8.727	ng	96
60) Diethylphthalate	10.339	149	122014	8.989	ng	99
61) 4-Chlorophenyl-phenyle...	10.463	204	62740	9.203	ng	100
62) 4-Nitroaniline	10.475	138	25720	8.105	ng	97
63) Azobenzene	10.622	77	127133	8.887	ng	99
65) 4,6-Dinitro-2-methylph...	10.504	198	10204	5.384	ng	97
66) n-Nitrosodiphenylamine	10.580	169	113468	9.504	ng	98
67) 4-Bromophenyl-phenylether	10.957	248	35492	8.919	ng	97
68) Hexachlorobenzene	11.016	284	40577	8.907	ng	99
69) Atrazine	11.104	200	31154	11.310	ng	98
70) Pentachlorophenol	11.210	266	16409	6.236	ng	95
71) Phenanthrene	11.433	178	182554	9.528	ng	99
72) Anthracene	11.480	178	181582	9.701	ng	98
73) Carbazole	11.639	167	161783	9.248	ng	98
74) Di-n-butylphthalate	11.969	149	162398	9.066	ng	99
75) Fluoranthene	12.621	202	179004	9.319	ng	98
77) Benzidine	12.745	184	44242	10.666	ng	96
78) Pyrene	12.851	202	180679	9.080	ng	99
80) Butylbenzylphthalate	13.469	149	37916	8.094	ng	98
81) Benzo(a)anthracene	14.033	228	115765	8.801	ng	98
82) 3,3'-Dichlorobenzidine	13.998	252	31012	8.878	ng	99
83) Chrysene	14.068	228	112910	9.197	ng	99
84) Bis(2-ethylhexyl)phtha...	14.033	149	42301	7.693	ng	99
85) Di-n-octyl phthalate	14.645	149	63771	6.058	ng	# 99
87) Indeno(1,2,3-cd)pyrene	16.992	276	83790	7.210	ng	99
88) Benzo(b)fluoranthene	15.086	252	87929	8.207	ng	99
89) Benzo(k)fluoranthene	15.115	252	97147	10.113	ng	98
90) Benzo(a)pyrene	15.457	252	81872	9.000	ng	98
91) Dibenzo(a,h)anthracene	17.015	278	64495	6.734	ng	99
92) Benzo(g,h,i)perylene	17.439	276	66519	6.729	ng	97

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

