

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081524\
 Data File : BP021509.D
 Acq On : 15 Aug 2024 11:18
 Operator : RC/JU
 Sample : P3390-03
 Misc : MDL-MDL-SOIL-4ng
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MDL-SOIL-03-QT3-2024

Quant Time: Aug 15 11:56:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 00:00:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	444773	20.000	ng	0.00
21) Naphthalene-d8	10.587	136	2011295	20.000	ng	-0.01
39) Acenaphthene-d10	14.439	164	1445135	20.000	ng	-0.02
64) Phenanthrene-d10	17.233	188	3303546	20.000	ng	-0.02
76) Chrysene-d12	21.698	240	3096346	20.000	ng	0.00
86) Perylene-d12	25.098	264	2978689	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	3878515	129.483	ng	0.00
7) Phenol-d6	6.963	99	5689916	130.217	ng	0.00
23) Nitrobenzene-d5	8.946	82	3843359	98.630	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	2331991	144.990	ng	-0.02
45) 2-Fluorobiphenyl	13.057	172	7970155	84.170	ng	0.00
79) Terphenyl-d14	19.969	244	12072471	75.746	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.317	88	60041	4.227	ng	99
3) Pyridine	3.711	79	143062	3.608	ng	98
4) n-Nitrosodimethylamine	3.617	42	52496	4.405	ng	# 91
6) Aniline	7.128	93	215030	4.922	ng	99
8) 2-Chlorophenol	7.369	128	120282	4.345	ng	99
9) Benzaldehyde	6.946	77	157870	5.643	ng	98
10) Phenol	6.987	94	189271	4.183	ng	95
11) bis(2-Chloroethyl)ether	7.228	93	166182	4.340	ng	98
12) 1,3-Dichlorobenzene	7.693	146	135633	4.125	ng	97
13) 1,4-Dichlorobenzene	7.840	146	137548	4.143	ng	99
14) 1,2-Dichlorobenzene	8.157	146	133464	4.171	ng	97
15) Benzyl Alcohol	8.028	79	137737	4.393	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.316	45	167093	4.739	ng	96
17) 2-Methylphenol	8.228	107	116128	4.060	ng	98
18) Hexachloroethane	8.887	117	55633	4.179	ng	98
19) n-Nitroso-di-n-propyla...	8.599	70	138328	4.583	ng	99
20) 3+4-Methylphenols	8.552	107	158041	4.097	ng	97
22) Acetophenone	8.610	105	264968	4.288	ng	# 99
24) Nitrobenzene	8.987	77	195237	4.613	ng	100
25) Isophorone	9.510	82	392641	4.265	ng	99
26) 2-Nitrophenol	9.699	139	31850	5.215	ng	94
27) 2,4-Dimethylphenol	9.757	122	85430	3.760	ng	98
28) bis(2-Chloroethoxy)met...	9.993	93	242002	4.307	ng	97
29) 2,4-Dichlorophenol	10.228	162	100866	3.758	ng	96
30) 1,2,4-Trichlorobenzene	10.451	180	137478	3.995	ng	97
31) Naphthalene	10.640	128	432576	4.117	ng	100
32) Benzoic acid	9.804	122	24885	5.622	ng	94
33) 4-Chloroaniline	10.740	127	175678	4.451	ng	98
34) Hexachlorobutadiene	10.934	225	84398	3.895	ng	94
35) Caprolactam	11.475	113	45756	4.095	ng	94
36) 4-Chloro-3-methylphenol	11.851	107	146797	3.942	ng	96
37) 2-Methylnaphthalene	12.251	142	297157	4.144	ng	98
38) 1-Methylnaphthalene	12.469	142	286232	4.047	ng	100
40) 1,2,4,5-Tetrachloroben...	12.616	216	170353	4.023	ng	99
41) Hexachlorocyclopentadiene	12.610	237	48723	3.631	ng	98
43) 2,4,6-Trichlorophenol	12.851	196	82308	3.499	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.916	196	94548	3.618	ng	97
46) 1,1'-Biphenyl	13.263	154	430469	4.210	ng	100
47) 2-Chloronaphthalene	13.304	162	318437	4.000	ng	98
48) 2-Nitroaniline	13.492	65	70086	7.952	ng	96
49) Acenaphthylene	14.157	152	519429	4.439	ng	99
50) Dimethylphthalate	13.887	163	425703	4.399	ng	100
51) 2,6-Dinitrotoluene	13.998	165	63940	6.988	ng	94
52) Acenaphthene	14.504	154	313380	4.082	ng	99
53) 3-Nitroaniline	14.328	138	65445	7.097	ng #	89
54) 2,4-Dinitrophenol	14.528	184	14845	6.395	ng #	69
55) Dibenzofuran	14.839	168	519511	4.398	ng	99
56) 4-Nitrophenol	14.628	139	47094	7.826	ng	91
57) 2,4-Dinitrotoluene	14.787	165	70860	7.577	ng	91
58) Fluorene	15.498	166	423248	4.362	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.069	232	89654	4.042	ng #	98
60) Diethylphthalate	15.275	149	436537	4.396	ng	99
61) 4-Chlorophenyl-phenylether	15.498	204	218063	4.348	ng	99
62) 4-Nitroaniline	15.498	138	69835	6.258	ng	92
63) Azobenzene	15.798	77	550767	4.512	ng	99
65) 4,6-Dinitro-2-methylph...	15.569	198	23385	7.584	ng	94
66) n-Nitrosodiphenylamine	15.716	169	364586	4.118	ng	100
67) 4-Bromophenyl-phenylether	16.410	248	128341	3.671	ng	99
68) Hexachlorobenzene	16.528	284	161936	3.919	ng	94
69) Atrazine	16.686	200	126177	4.151	ng	99
70) Pentachlorophenol	16.875	266	66342	2.942	ng	95
71) Phenanthrene	17.281	178	683628	4.303	ng	100
72) Anthracene	17.375	178	648787	4.164	ng	99
73) Carbazole	17.651	167	608876	4.055	ng	99
74) Di-n-butylphthalate	18.251	149	708923	3.844	ng	99
75) Fluoranthene	19.363	202	773721	3.969	ng	99
77) Benzidine	19.557	184	211373	6.819	ng	99
78) Pyrene	19.739	202	836851	4.135	ng	100
80) Butylbenzylphthalate	20.704	149	208104	6.108	ng	94
81) Benzo(a)anthracene	21.674	228	800948	4.121	ng	100
82) 3,3'-Dichlorobenzidine	21.598	252	216395	3.472	ng	98
83) Chrysene	21.745	228	759519	4.139	ng	99
84) Bis(2-ethylhexyl)phtha...	21.639	149	377936	3.290	ng	99
85) Di-n-octyl phthalate	22.951	149	444984	6.819	ng	97
87) Indeno(1,2,3-cd)pyrene	28.962	276	628413	3.255	ng	97
88) Benzo(b)fluoranthene	24.021	252	684912	4.004	ng	98
89) Benzo(k)fluoranthene	24.092	252	673993	4.039	ng	98
90) Benzo(a)pyrene	24.927	252	557245	3.754	ng	98
91) Dibenzo(a,h)anthracene	29.050	278	525517	3.276	ng	99
92) Benzo(g,h,i)perylene	30.145	276	505802	3.146	ng	97

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

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