

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092024\
 Data File : BP021963.D
 Acq On : 20 Sep 2024 11:49
 Operator : RC/JU
 Sample : P2403-02
 Misc : LOQ--SOIL 5 ppm
 ALS Vial : 6 Sample Multiplier: 1

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

Quant Time: Sep 20 12:33:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 01:22:00 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/20/2024
 Supervised By :mohammad ahmed 09/24/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	75132	20.000	ng	0.00
21) Naphthalene-d8	10.475	136	282360	20.000	ng	0.00
39) Acenaphthene-d10	14.322	164	208918	20.000	ng	0.00
64) Phenanthrene-d10	17.110	188	508713	20.000	ng	-0.03
76) Chrysene-d12	21.539	240	626667	20.000	ng	-0.02
86) Perylene-d12	24.833	264	750639	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	468864	117.277	ng	0.00
7) Phenol-d6	6.893	99	603181	116.219	ng	0.00
23) Nitrobenzene-d5	8.852	82	456145	85.480	ng	0.00
42) 2,4,6-Tribromophenol	15.834	330	482980	124.701	ng	-0.02
45) 2-Fluorobiphenyl	12.934	172	1227826	85.792	ng	0.00
79) Terphenyl-d14	19.845	244	2734781	81.977	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.258	88	8001	4.820	ng	94
3) Pyridine	3.664	79	18607	4.203	ng	98
4) n-Nitrosodimethylamine	3.564	42	11323	4.848	ng	# 86
6) Aniline	7.046	93	26796	5.875	ng	93
8) 2-Chlorophenol	7.281	128	21101	4.936	ng	91
9) Benzaldehyde	6.864	77	21032	7.750	ng	95
10) Phenol	6.917	94	25009	4.852	ng	96
11) bis(2-Chloroethyl)ether	7.140	93	19043	4.904	ng	89
12) 1,3-Dichlorobenzene	7.605	146	25901	4.799	ng	95
13) 1,4-Dichlorobenzene	7.746	146	26037	4.757	ng	97
14) 1,2-Dichlorobenzene	8.058	146	25043	4.767	ng	98
15) Benzyl Alcohol	7.952	79	18191	4.570	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.222	45	21268	5.135	ng	98
17) 2-Methylphenol	8.146	107	16130	4.632	ng	97
18) Hexachloroethane	8.775	117	9346	4.644	ng	97
19) n-Nitroso-di-n-propyla...	8.505	70	15913	4.933	ng	92
20) 3+4-Methylphenols	8.469	107	21218	4.316	ng	96
22) Acetophenone	8.522	105	34001	4.936	ng	97
24) Nitrobenzene	8.893	77	24987	4.611	ng	99
25) Isophorone	9.416	82	41227	4.686	ng	97
26) 2-Nitrophenol	9.599	139	9640	4.068	ng	88
27) 2,4-Dimethylphenol	9.652	122	12404	4.521	ng	94
28) bis(2-Chloroethoxy)met...	9.887	93	24467	4.638	ng	97
29) 2,4-Dichlorophenol	10.128	162	19953	4.285	ng	96
30) 1,2,4-Trichlorobenzene	10.340	180	27073	4.667	ng	96
31) Naphthalene	10.528	128	66378	4.711	ng	97
32) Benzoic acid	9.728	122	10392m	3.240	ng	
33) 4-Chloroaniline	10.628	127	25199	4.958	ng	94
34) Hexachlorobutadiene	10.810	225	20567	4.532	ng	98
35) Caprolactam	11.399	113	5338	3.852	ng	# 89
36) 4-Chloro-3-methylphenol	11.757	107	20240	4.051	ng	# 91
37) 2-Methylnaphthalene	12.128	142	48768	4.731	ng	98
38) 1-Methylnaphthalene	12.352	142	43855	4.309	ng	100
40) 1,2,4,5-Tetrachloroben...	12.504	216	34491	4.719	ng	99
41) Hexachlorocyclopentadiene	12.487	237	15090	5.556	ng	97
43) 2,4,6-Trichlorophenol	12.746	196	18405	4.102	ng	93

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44) 2,4,5-Trichlorophenol	12.816	196	19660	3.906	ng	97
46) 1,1'-Biphenyl	13.146	154	73761	5.178	ng	97
47) 2-Chloronaphthalene	13.193	162	53911	4.681	ng	98
48) 2-Nitroaniline	13.393	65	12748	4.053	ng	99
49) Acenaphthylene	14.046	152	79400	4.885	ng	98
50) Dimethylphthalate	13.775	163	73562	4.855	ng	99
51) 2,6-Dinitrotoluene	13.887	165	13543	4.011	ng	97
52) Acenaphthene	14.381	154	48002	4.538	ng	98
53) 3-Nitroaniline	14.228	138	11794	3.919	ng	99
54) 2,4-Dinitrophenol	14.440	184	5921	2.993	ng #	77
55) Dibenzofuran	14.722	168	87727	4.934	ng	99
56) 4-Nitrophenol	14.551	139	8334	3.182	ng	85
57) 2,4-Dinitrotoluene	14.687	165	18641	3.905	ng #	93
58) Fluorene	15.381	166	69789	4.700	ng	93
59) 2,3,4,6-Tetrachlorophenol	14.957	232	21192	4.363	ng	99
60) Diethylphthalate	15.151	149	72756	4.725	ng	99
61) 4-Chlorophenyl-phenyle...	15.375	204	39801	4.604	ng	97
62) 4-Nitroaniline	15.393	138	12893	3.896	ng	97
63) Azobenzene	15.675	77	57578	4.519	ng	98
65) 4,6-Dinitro-2-methylph...	15.463	198	10330	3.503	ng	95
66) n-Nitrosodiphenylamine	15.593	169	57832	4.566	ng	98
67) 4-Bromophenyl-phenylether	16.287	248	27278	4.600	ng	98
68) Hexachlorobenzene	16.404	284	34530	4.670	ng	93
69) Atrazine	16.563	200	25342	5.584	ng	97
70) Pentachlorophenol	16.757	266	17749	3.568	ng	96
71) Phenanthrene	17.157	178	120610	4.939	ng	97
72) Anthracene	17.251	178	112788	4.771	ng	98
73) Carbazole	17.534	167	101661	4.453	ng	99
74) Di-n-butylphthalate	18.122	149	112604	4.199	ng	99
75) Fluoranthene	19.239	202	152606	4.477	ng	99
77) Benzidine	19.439	184	63355	8.273	ng	96
78) Pyrene	19.616	202	162904	4.297	ng	98
80) Butylbenzylphthalate	20.575	149	55953	4.138	ng	93
81) Benzo(a)anthracene	21.522	228	192887	4.790	ng	99
82) 3,3'-Dichlorobenzidine	21.445	252	64301	4.401	ng	100
83) Chrysene	21.592	228	184677	4.871	ng	98
84) Bis(2-ethylhexyl)phtha...	21.469	149	84160	4.241	ng	99
85) Di-n-octyl phthalate	22.710	149	136346	4.045	ng	99
87) Indeno(1,2,3-cd)pyrene	28.551	276	240368	4.815	ng #	88
88) Benzo(b)fluoranthene	23.786	252	199947m	4.504	ng	
89) Benzo(k)fluoranthene	23.845	252	203228	4.837	ng	98
90) Benzo(a)pyrene	24.669	252	179516	4.694	ng	97
91) Dibenzo(a,h)anthracene	28.633	278	197614	4.796	ng #	97
92) Benzo(g,h,i)perylene	29.698	276	188242	4.460	ng #	97

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

