

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042924\
 Data File : BF137748.D
 Acq On : 29 Apr 2024 11:16
 Operator : CG\JU
 Sample : PB160543BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB160543BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/30/2024
 Supervised By :mohammad ahmed 05/03/2024

Quant Time: Apr 29 11:52:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 03:57:12 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.063	152	345310	20.000	ng	0.00
21) Naphthalene-d8	8.351	136	1392838	20.000	ng	0.00
39) Acenaphthene-d10	10.110	164	735841	20.000	ng	0.00
64) Phenanthrene-d10	11.610	188	1244526	20.000	ng	0.00
76) Chrysene-d12	14.257	240	742027	20.000	ng	0.00
86) Perylene-d12	15.827	264	622181	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.704	112	2750350	120.989	ng	0.02
7) Phenol-d6	6.687	99	3469675	120.194	ng	0.01
23) Nitrobenzene-d5	7.634	82	2214565	84.401	ng	0.00
42) 2,4,6-Tribromophenol	10.910	330	1021310	137.547	ng	0.00
45) 2-Fluorobiphenyl	9.428	172	4067849	82.602	ng	0.00
79) Terphenyl-d14	13.198	244	4171681	90.490	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.116	88	346319	34.023	ng	Qvalue 97
3) Pyridine	3.840	79	812861	29.977	ng	97
4) n-Nitrosodimethylamine	3.810	42	505247	39.925	ng	92
6) Aniline	6.728	93	585283	16.290	ng	98
8) 2-Chlorophenol	6.851	128	1121660	47.269	ng	100
10) Phenol	6.698	94	1275046	44.213	ng	97
11) bis(2-Chloroethyl)ether	6.798	93	1054035	44.810	ng	98
12) 1,3-Dichlorobenzene	7.004	146	1131440	45.467	ng	98
13) 1,4-Dichlorobenzene	7.081	146	1155177	46.108	ng	99
14) 1,2-Dichlorobenzene	7.234	146	1118410	47.958	ng	99
15) Benzyl Alcohol	7.204	79	940801	43.566	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.328	45	1529995	42.294	ng	98
17) 2-Methylphenol	7.304	107	896754	42.535	ng	100
18) Hexachloroethane	7.581	117	413133	46.592	ng	96
19) n-Nitroso-di-n-propyla...	7.475	70	763234	40.887	ng	99
20) 3+4-Methylphenols	7.457	107	1121907	41.201	ng	# 89
22) Acetophenone	7.475	105	1486035	43.621	ng	100
24) Nitrobenzene	7.651	77	1150252	43.391	ng	99
25) Isophorone	7.887	82	2137746	43.868	ng	99
26) 2-Nitrophenol	7.963	139	601135	51.674	ng	97
27) 2,4-Dimethylphenol	7.987	122	908086	38.691	ng	98
28) bis(2-Chloroethoxy)met...	8.092	93	1294169	43.480	ng	99
29) 2,4-Dichlorophenol	8.198	162	952372	46.301	ng	97
30) 1,2,4-Trichlorobenzene	8.287	180	988430	47.145	ng	98
31) Naphthalene	8.375	128	3127077	43.996	ng	100
32) Benzoic acid	8.116	122	765362	51.502	ng	98
33) 4-Chloroaniline	8.410	127	437328	14.537	ng	98
34) Hexachlorobutadiene	8.481	225	586221	46.641	ng	99
35) Caprolactam	8.798	113	264166m	39.723	ng	
36) 4-Chloro-3-methylphenol	8.892	107	976692	44.769	ng	95
37) 2-Methylnaphthalene	9.063	142	2031904	41.651	ng	99
38) 1-Methylnaphthalene	9.163	142	1936736	42.462	ng	100
40) 1,2,4,5-Tetrachloroben...	9.228	216	981799	48.668	ng	100
41) Hexachlorocyclopentadiene	9.216	237	1006721	89.064	ng	100
43) 2,4,6-Trichlorophenol	9.333	196	697595	49.293	ng	97
44) 2,4,5-Trichlorophenol	9.392	196	702138	43.291	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.528	154	2449398	45.757	ng	100
47) 2-Chloronaphthalene	9.557	162	1949744	47.250	ng	98
48) 2-Nitroaniline	9.651	65	590262	46.942	ng	92
49) Acenaphthylene	9.975	152	2996889	46.571	ng	99
50) Dimethylphthalate	9.828	163	2311688	45.090	ng	99
51) 2,6-Dinitrotoluene	9.892	165	527090	48.796	ng	87
52) Acenaphthene	10.151	154	2057018	49.367	ng	99
53) 3-Nitroaniline	10.063	138	368383	30.347	ng	99
54) 2,4-Dinitrophenol	10.169	184	610923	104.600	ng	# 45
55) Dibenzofuran	10.322	168	2713092	44.419	ng	99
56) 4-Nitrophenol	10.216	139	890233	94.302	ng	98
57) 2,4-Dinitrotoluene	10.298	165	700138	51.413	ng	97
58) Fluorene	10.669	166	2102413	44.052	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.433	232	590639	47.030	ng	97
60) Diethylphthalate	10.528	149	2135424	43.306	ng	99
61) 4-Chlorophenyl-phenyle...	10.651	204	1056113	44.785	ng	97
62) 4-Nitroaniline	10.686	138	521434	45.098	ng	99
63) Azobenzene	10.816	77	2086175	41.286	ng	94
65) 4,6-Dinitro-2-methylph...	10.710	198	386838	54.212	ng	# 75
66) n-Nitrosodiphenylamine	10.775	169	1884927	44.719	ng	99
67) 4-Bromophenyl-phenylether	11.145	248	669784	46.445	ng	98
68) Hexachlorobenzene	11.210	284	705958	50.250	ng	100
69) Atrazine	11.292	200	427043	36.650	ng	99
70) Pentachlorophenol	11.404	266	909043	86.017	ng	99
71) Phenanthrene	11.633	178	2900319	44.025	ng	99
72) Anthracene	11.686	178	2925414	43.365	ng	100
73) Carbazole	11.839	167	2647660	43.307	ng	99
74) Di-n-butylphthalate	12.157	149	3111744	41.753	ng	99
75) Fluoranthene	12.827	202	2892363	41.769	ng	99
77) Benzidine	12.939	184	179161	10.337	ng	99
78) Pyrene	13.057	202	2928589	48.085	ng	100
80) Butylbenzylphthalate	13.663	149	1096869	50.785	ng	94
81) Benzo(a)anthracene	14.251	228	2387301	46.127	ng	99
82) 3,3'-Dichlorobenzidine	14.204	252	560526	35.004	ng	98
83) Chrysene	14.286	228	2165136	44.773	ng	99
84) Bis(2-ethylhexyl)phtha...	14.216	149	1528884	46.873	ng	100
85) Di-n-octyl phthalate	14.845	149	2291613	46.772	ng	97
87) Indeno(1,2,3-cd)pyrene	17.480	276	2142754	62.112	ng	99
88) Benzo(b)fluoranthene	15.357	252	2047704	51.432	ng	99
89) Benzo(k)fluoranthene	15.392	252	1886942	48.214	ng	99
90) Benzo(a)pyrene	15.768	252	1743118	48.805	ng	97
91) Dibenzo(a,h)anthracene	17.504	278	1782962	62.279	ng	99
92) Benzo(g,h,i)perylene	17.986	276	1717942	61.711	ng	98

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

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