

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042622\
 Data File : BF127906.D
 Acq On : 26 Apr 2022 11:18
 Operator : CG\JU
 Sample : PB144358BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB144358BS

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

04/27/2022
 Supervised By :Jagrut
 Upadhyay

Quant Time: Apr 27 01:46:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 00:43:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.946	152	186544	20.000	ng	0.00
21) Naphthalene-d8	8.228	136	727957	20.000	ng	0.00
39) Acenaphthene-d10	9.992	164	398458	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	671386	20.000	ng	0.01
76) Chrysene-d12	14.133	240	406032	20.000	ng	0.00
86) Perylene-d12	15.645	264	344855	20.000	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	1016680	86.808	ng	0.02
7) Phenol-d6	6.575	99	1325466	87.220	ng	0.00
23) Nitrobenzene-d5	7.516	82	1369229	88.449	ng	0.00
42) 2,4,6-Tribromophenol	10.786	330	347071	86.526	ng	0.00
45) 2-Fluorobiphenyl	9.310	172	2082337	88.410	ng	0.00
79) Terphenyl-d14	13.069	244	2033670	91.586	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.952	88	199942	35.781	ng	98
3) Pyridine	3.687	79	553527	38.660	ng	99
4) n-Nitrosodimethylamine	3.663	42	327786	47.296	ng	100
6) Aniline	6.610	93	689185	38.170	ng	99
8) 2-Chlorophenol	6.734	128	589953	48.533	ng	99
9) Benzaldehyde	6.498	77	401203	42.675	ng	93
10) Phenol	6.587	94	717390m	46.805	ng	
11) bis(2-Chloroethyl)ether	6.681	93	608194	47.835	ng	98
12) 1,3-Dichlorobenzene	6.887	146	621468	44.415	ng	98
13) 1,4-Dichlorobenzene	6.963	146	626036	44.824	ng	98
14) 1,2-Dichlorobenzene	7.116	146	591154	45.722	ng	99
15) Benzyl Alcohol	7.093	79	503268	48.698	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.216	45	919762	47.331	ng	97
17) 2-Methylphenol	7.198	107	527204	50.267	ng	95
18) Hexachloroethane	7.457	117	244903	47.697	ng	98
19) n-Nitroso-di-n-propyla...	7.363	70	443040	47.625	ng	99
20) 3+4-Methylphenols	7.351	107	547082	43.180	ng	92
22) Acetophenone	7.357	105	782279	44.090	ng	# 91
24) Nitrobenzene	7.534	77	690770	46.307	ng	98
25) Isophorone	7.769	82	1303103	50.014	ng	99
26) 2-Nitrophenol	7.845	139	313458	49.312	ng	97
27) 2,4-Dimethylphenol	7.881	122	554264	52.496	ng	100
28) bis(2-Chloroethoxy)met...	7.975	93	766845	45.925	ng	98
29) 2,4-Dichlorophenol	8.087	162	490448	44.340	ng	98
30) 1,2,4-Trichlorobenzene	8.169	180	553086	44.020	ng	100
31) Naphthalene	8.251	128	1637872	44.176	ng	100
32) Benzoic acid	8.016	122	301190	39.799	ng	98
33) 4-Chloroaniline	8.298	127	350033	23.861	ng	99
34) Hexachlorobutadiene	8.363	225	358484	45.297	ng	98
35) Caprolactam	8.687	113	148859m	41.692	ng	
36) 4-Chloro-3-methylphenol	8.787	107	534941	45.284	ng	98
37) 2-Methylnaphthalene	8.945	142	1101744	44.963	ng	99
38) 1-Methylnaphthalene	9.045	142	1039445	43.643	ng	99
40) 1,2,4,5-Tetrachloroben...	9.110	216	513828	43.750	ng	98
41) Hexachlorocyclopentadiene	9.092	237	719828	113.125	ng	99
43) 2,4,6-Trichlorophenol	9.222	196	352239	45.580	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.263	196	361944	43.084	ng	99
46) 1,1'-Biphenyl	9.410	154	1322732	44.815	ng	99
47) 2-Chloronaphthalene	9.434	162	975548	43.417	ng	98
48) 2-Nitroaniline	9.534	65	353240	46.476	ng	94
49) Acenaphthylene	9.857	152	1557016	45.301	ng	100
50) Dimethylphthalate	9.716	163	1166758	43.662	ng	99
51) 2,6-Dinitrotoluene	9.775	165	266905	46.131	ng	94
52) Acenaphthene	10.028	154	1107424m	47.925	ng	
53) 3-Nitroaniline	9.945	138	170552	26.546	ng	98
54) 2,4-Dinitrophenol	10.057	184	279764	91.815	ng	# 85
55) Dibenzofuran	10.198	168	1375160	41.248	ng	98
56) 4-Nitrophenol	10.116	139	376727	76.789	ng	95
57) 2,4-Dinitrotoluene	10.186	165	347237	45.308	ng	93
58) Fluorene	10.545	166	1036572	41.668	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.316	232	300193	42.494	ng	100
60) Diethylphthalate	10.410	149	1114103	42.516	ng	100
61) 4-Chlorophenyl-phenyle...	10.534	204	517170	41.835	ng	98
62) 4-Nitroaniline	10.569	138	260673	41.500	ng	98
63) Azobenzene	10.692	77	1206044	41.827	ng	99
65) 4,6-Dinitro-2-methylph...	10.592	198	181686	45.543	ng	95
66) n-Nitrosodiphenylamine	10.651	169	922827	44.174	ng	99
67) 4-Bromophenyl-phenylether	11.022	248	331372	44.523	ng	99
68) Hexachlorobenzene	11.092	284	360529	45.977	ng	# 90
69) Atrazine	11.181	200	330605	51.179	ng	98
70) Pentachlorophenol	11.286	266	386052	83.684	ng	98
71) Phenanthrene	11.510	178	1499976	42.681	ng	100
72) Anthracene	11.563	178	1518672	43.380	ng	99
73) Carbazole	11.716	167	1338521	39.671	ng	99
74) Di-n-butylphthalate	12.033	149	1649084	43.465	ng	99
75) Fluoranthene	12.704	202	1521533	41.136	ng	97
77) Benzidine	12.822	184	707427	92.080	ng	98
78) Pyrene	12.933	202	1536569	46.880	ng	100
80) Butylbenzylphthalate	13.539	149	643933	50.123	ng	99
81) Benzo(a)anthracene	14.121	228	1228231	45.384	ng	98
82) 3,3'-Dichlorobenzidine	14.080	252	253949	29.553	ng	100
83) Chrysene	14.163	228	1127039	43.605	ng	100
84) Bis(2-ethylhexyl)phtha...	14.092	149	883394	51.852	ng	99
85) Di-n-octyl phthalate	14.716	149	1357359	51.411	ng	100
87) Indeno(1,2,3-cd)pyrene	17.204	276	1044066	45.358	ng	99
88) Benzo(b)fluoranthene	15.198	252	1016561	46.933	ng	99
89) Benzo(k)fluoranthene	15.227	252	1007396	46.164	ng	97
90) Benzo(a)pyrene	15.586	252	968090	53.882	ng	99
91) Dibenzo(a,h)anthracene	17.227	278	894783	48.470	ng	98
92) Benzo(g,h,i)perylene	17.680	276	910148	46.783	ng	99

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

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