

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111622\
 Data File : BF131260.D
 Acq On : 16 Nov 2022 12:23
 Operator : CG\JU
 Sample : PB148898BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB148898BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/17/2022
 Supervised By : mohammad ahmed 11/17/2022

Quant Time: Nov 16 22:13:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 06:35:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.975	152	95509	20.000	ng	0.00
21) Naphthalene-d8	8.269	136	347686	20.000	ng	0.00
39) Acenaphthene-d10	10.033	164	197310	20.000	ng	0.00
64) Phenanthrene-d10	11.527	188	366898	20.000	ng	0.00
76) Chrysene-d12	14.186	240	252563	20.000	ng	0.00
86) Perylene-d12	15.727	264	191277	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.592	112	733138	135.925	ng	0.02
7) Phenol-d6	6.592	99	918507	134.452	ng	0.00
23) Nitrobenzene-d5	7.545	82	591766	95.534	ng	0.00
42) 2,4,6-Tribromophenol	10.827	330	264406	140.289	ng	0.00
45) 2-Fluorobiphenyl	9.351	172	1106026	82.032	ng	0.00
79) Terphenyl-d14	13.127	244	1358916	92.398	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.887	88	91140	36.882	ng	98
3) Pyridine	3.628	79	240427	34.422	ng	99
4) n-Nitrosodimethylamine	3.604	42	179714	44.144	ng	96
6) Aniline	6.634	93	247238m	28.853	ng	
8) 2-Chlorophenol	6.757	128	269894	44.514	ng	99
9) Benzaldehyde	6.522	77	48390	9.765	ng	95
10) Phenol	6.604	94	321923	42.124	ng	95
11) bis(2-Chloroethyl)ether	6.710	93	252994	42.280	ng	99
12) 1,3-Dichlorobenzene	6.916	146	286887	41.824	ng	99
13) 1,4-Dichlorobenzene	6.992	146	288387	41.572	ng	98
14) 1,2-Dichlorobenzene	7.145	146	281043	43.833	ng	99
15) Benzyl Alcohol	7.116	79	216393m	36.467	ng	
16) 2,2'-oxybis(1-Chloropr...	7.245	45	488372	43.331	ng	100
17) 2-Methylphenol	7.222	107	225436	43.553	ng	98
18) Hexachloroethane	7.492	117	109721	43.507	ng	93
19) n-Nitroso-di-n-propyla...	7.392	70	201333	42.294	ng	97
20) 3+4-Methylphenols	7.375	107	281046	41.976	ng	# 88
22) Acetophenone	7.386	105	388003	42.811	ng	# 99
24) Nitrobenzene	7.563	77	304446	45.528	ng	98
25) Isophorone	7.798	82	542768	43.861	ng	98
26) 2-Nitrophenol	7.881	139	121688	45.647	ng	97
27) 2,4-Dimethylphenol	7.910	122	276784	54.644	ng	98
28) bis(2-Chloroethoxy)met...	8.010	93	310943	44.567	ng	99
29) 2,4-Dichlorophenol	8.116	162	230507	42.081	ng	99
30) 1,2,4-Trichlorobenzene	8.204	180	264080	41.119	ng	99
31) Naphthalene	8.286	128	746946	39.653	ng	98
32) Benzoic acid	8.028	122	142380m	40.080	ng	
33) 4-Chloroaniline	8.333	127	93788	12.648	ng	95
34) Hexachlorobutadiene	8.404	225	168645	41.490	ng	98
35) Caprolactam	8.704	113	73405m	45.346	ng	
36) 4-Chloro-3-methylphenol	8.816	107	249172	43.215	ng	95
37) 2-Methylnaphthalene	8.986	142	505084	40.260	ng	100
38) 1-Methylnaphthalene	9.086	142	477413	39.245	ng	100
40) 1,2,4,5-Tetrachloroben...	9.145	216	261117	41.102	ng	99
41) Hexachlorocyclopentadiene	9.133	237	334084	102.602	ng	97
43) 2,4,6-Trichlorophenol	9.257	196	150936	36.340	ng	99

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44) 2,4,5-Trichlorophenol	9.298	196	198893	44.624	ng	99
46) 1,1'-Biphenyl	9.451	154	634132	41.807	ng	99
47) 2-Chloronaphthalene	9.475	162	501204	40.790	ng	99
48) 2-Nitroaniline	9.569	65	180296	45.919	ng	97
49) Acenaphthylene	9.898	152	770644	41.072	ng	99
50) Dimethylphthalate	9.751	163	586890	41.538	ng	99
51) 2,6-Dinitrotoluene	9.816	165	126755	45.876	ng	96
52) Acenaphthene	10.069	154	519010	44.314	ng	99
53) 3-Nitroaniline	9.980	138	67545	24.254	ng	94
54) 2,4-Dinitrophenol	10.092	184	108626	84.621	ng	97
55) Dibenzofuran	10.245	168	678292	40.422	ng	99
56) 4-Nitrophenol	10.139	139	181744	81.755	ng	96
57) 2,4-Dinitrotoluene	10.222	165	160002	46.634	ng	94
58) Fluorene	10.586	166	537587	40.713	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.351	232	163954	43.422	ng	99
60) Diethylphthalate	10.457	149	561502	42.043	ng	99
61) 4-Chlorophenyl-phenyle...	10.580	204	270807	41.477	ng	95
62) 4-Nitroaniline	10.610	138	119845	41.386	ng	97
63) Azobenzene	10.739	77	572055	40.911	ng	99
65) 4,6-Dinitro-2-methylph...	10.627	198	64663	40.651	ng	88
66) n-Nitrosodiphenylamine	10.698	169	474577	40.713	ng	99
67) 4-Bromophenyl-phenylether	11.074	248	181967	42.240	ng	# 91
68) Hexachlorobenzene	11.133	284	195299	41.832	ng	96
69) Atrazine	11.227	200	162557	42.545	ng	97
70) Pentachlorophenol	11.327	266	218257m	80.096	ng	
71) Phenanthrene	11.557	178	809775	39.992	ng	99
72) Anthracene	11.610	178	835581	42.243	ng	99
73) Carbazole	11.763	167	736253	41.583	ng	99
74) Di-n-butylphthalate	12.092	149	860411	41.755	ng	100
75) Fluoranthene	12.751	202	916521	41.042	ng	99
77) Benzidine	12.874	184	187141	34.577	ng	98
78) Pyrene	12.986	202	935783	42.982	ng	99
80) Butylbenzylphthalate	13.604	149	345600	46.840	ng	99
81) Benzo(a)anthracene	14.180	228	728159	41.484	ng	99
82) 3,3'-Dichlorobenzidine	14.139	252	164402	32.009	ng	98
83) Chrysene	14.215	228	685376	40.082	ng	99
84) Bis(2-ethylhexyl)phtha...	14.162	149	434718	46.847	ng	99
85) Di-n-octyl phthalate	14.792	149	598938	45.674	ng	99
87) Indeno(1,2,3-cd)pyrene	17.321	276	550707	43.581	ng	99
88) Benzo(b)fluoranthene	15.274	252	578520	43.319	ng	99
89) Benzo(k)fluoranthene	15.304	252	570999	43.130	ng	99
90) Benzo(a)pyrene	15.668	252	488163	45.728	ng	99
91) Dibenzo(a,h)anthracene	17.345	278	472400	45.566	ng	99
92) Benzo(g,h,i)perylene	17.803	276	450347	39.500	ng	98

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

