

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070623\
 Data File : BF134103.D
 Acq On : 06 Jul 2023 18:15
 Operator : CG\JU
 Sample : PB153945BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB153945BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/07/2023
 Supervised By :mohammad ahmed 07/07/2023

Quant Time: Jul 07 01:30:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 30 18:35:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.846	152	108951	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	421083	20.000	ng	0.00
39) Acenaphthene-d10	9.887	164	212403	20.000	ng	0.00
64) Phenanthrene-d10	11.381	188	413777	20.000	ng	0.00
76) Chrysene-d12	14.045	240	255979	20.000	ng	0.00
86) Perylene-d12	15.539	264	204526	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	765665	117.556	ng	0.00
7) Phenol-d6	6.493	99	908200	113.384	ng	0.00
23) Nitrobenzene-d5	7.416	82	625188	80.034	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	306772	115.193	ng	0.00
45) 2-Fluorobiphenyl	9.204	172	1145987	78.443	ng	0.00
79) Terphenyl-d14	12.975	244	1341782	84.362	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.740	88	112358	41.837	ng	98
3) Pyridine	3.510	79	246276	35.206	ng	97
4) n-Nitrosodimethylamine	3.475	42	180384	45.149	ng	98
6) Aniline	6.510	93	240521	24.676	ng	# 74
8) 2-Chlorophenol	6.634	128	301842	45.653	ng	99
9) Benzaldehyde	6.399	77	37739	6.235	ng	98
10) Phenol	6.504	94	357354	42.432	ng	89
11) bis(2-Chloroethyl)ether	6.587	93	274386	44.253	ng	97
12) 1,3-Dichlorobenzene	6.787	146	312758	44.370	ng	99
13) 1,4-Dichlorobenzene	6.863	146	326474	45.855	ng	99
14) 1,2-Dichlorobenzene	7.016	146	305069	45.752	ng	99
15) Benzyl Alcohol	6.993	79	272210	44.083	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.122	45	452671	41.267	ng	97
17) 2-Methylphenol	7.104	107	252015	42.565	ng	97
18) Hexachloroethane	7.351	117	122762	46.503	ng	94
19) n-Nitroso-di-n-propyla...	7.263	70	210181	41.377	ng	99
20) 3+4-Methylphenols	7.257	107	301051	41.913	ng	95
22) Acetophenone	7.257	105	426978	44.586	ng	97
24) Nitrobenzene	7.434	77	334315	42.352	ng	97
25) Isophorone	7.669	82	583609	41.108	ng	100
26) 2-Nitrophenol	7.745	139	164826	43.527	ng	94
27) 2,4-Dimethylphenol	7.781	122	257462	42.952	ng	95
28) bis(2-Chloroethoxy)met...	7.875	93	334051	40.637	ng	99
29) 2,4-Dichlorophenol	7.987	162	248771	41.853	ng	97
30) 1,2,4-Trichlorobenzene	8.069	180	268219	43.733	ng	98
31) Naphthalene	8.151	128	822667	40.446	ng	98
32) Benzoic acid	7.916	122	194143	43.223	ng	96
33) 4-Chloroaniline	8.198	127	85396	9.815	ng	97
34) Hexachlorobutadiene	8.263	225	172031	44.466	ng	99
35) Caprolactam	8.575	113	83656m	42.745	ng	
36) 4-Chloro-3-methylphenol	8.687	107	288255	43.169	ng	97
37) 2-Methylnaphthalene	8.840	142	559618	38.907	ng	100
38) 1-Methylnaphthalene	8.940	142	519145	38.825	ng	99
40) 1,2,4,5-Tetrachloroben...	9.010	216	271659	43.170	ng	99
41) Hexachlorocyclopentadiene	8.992	237	273301	91.545	ng	100
43) 2,4,6-Trichlorophenol	9.122	196	178904	41.763	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.169	196	193405	38.382	ng	99
46) 1,1'-Biphenyl	9.310	154	712441	41.814	ng	98
47) 2-Chloronaphthalene	9.334	162	519161	41.645	ng	95
48) 2-Nitroaniline	9.434	65	191688	42.191	ng	98
49) Acenaphthylene	9.751	152	860499	40.408	ng	99
50) Dimethylphthalate	9.610	163	623447	40.435	ng	100
51) 2,6-Dinitrotoluene	9.675	165	139297	41.946	ng	88
52) Acenaphthene	9.922	154	508463	41.148	ng	100
53) 3-Nitroaniline	9.845	138	87966	22.080	ng	99
54) 2,4-Dinitrophenol	9.957	184	166363	86.420	ng	87
55) Dibenzofuran	10.098	168	768773	39.809	ng	99
56) 4-Nitrophenol	10.016	139	224547	80.578	ng	91
57) 2,4-Dinitrotoluene	10.086	165	184716	42.118	ng	94
58) Fluorene	10.439	166	608302	40.488	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.216	232	161401m	40.607	ng	
60) Diethylphthalate	10.316	149	646378	40.239	ng	99
61) 4-Chlorophenyl-phenyle...	10.428	204	287053	39.646	ng	94
62) 4-Nitroaniline	10.469	138	150538	37.494	ng	94
63) Azobenzene	10.592	77	603173	39.696	ng	98
65) 4,6-Dinitro-2-methylph...	10.498	198	110035	48.777	ng	100
66) n-Nitrosodiphenylamine	10.551	169	534212	40.408	ng	98
67) 4-Bromophenyl-phenylether	10.922	248	179254	40.758	ng	92
68) Hexachlorobenzene	10.986	284	206543	44.232	ng	96
69) Atrazine	11.081	200	148891	40.716	ng	99
70) Pentachlorophenol	11.186	266	210031	75.225	ng	98
71) Phenanthrene	11.410	178	866854	41.615	ng	99
72) Anthracene	11.457	178	887117	40.945	ng	99
73) Carbazole	11.616	167	853763	43.052	ng	99
74) Di-n-butylphthalate	11.945	149	986990	41.852	ng	99
75) Fluoranthene	12.604	202	952199	42.284	ng	98
77) Benzidine	12.728	184	133801	21.968	ng	97
78) Pyrene	12.833	202	964181	40.474	ng	99
80) Butylbenzylphthalate	13.457	149	383617	42.937	ng	95
81) Benzo(a)anthracene	14.033	228	741486	41.768	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	167235	29.734	ng #	97
83) Chrysene	14.069	228	696785	40.524	ng	100
84) Bis(2-ethylhexyl)phtha...	14.016	149	443586	43.208	ng	99
85) Di-n-octyl phthalate	14.633	149	638279	42.312	ng	99
87) Indeno(1,2,3-cd)pyrene	17.092	276	659541	46.473	ng	98
88) Benzo(b)fluoranthene	15.098	252	600728	49.951	ng	100
89) Benzo(k)fluoranthene	15.133	252	564138	46.072	ng	98
90) Benzo(a)pyrene	15.480	252	501207	43.099	ng	98
91) Dibenzo(a,h)anthracene	17.110	278	541905	46.748	ng	97
92) Benzo(g,h,i)perylene	17.568	276	559566	46.810	ng	97

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

