

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112822\
 Data File : BF131422.D
 Acq On : 28 Nov 2022 10:38
 Operator : CG\JU
 Sample : PB149277BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB149277BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/02/2022
 Supervised By : mohammad ahmed 12/02/2022

Quant Time: Nov 29 00:19:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 00:18:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.951	152	86833	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	320955	20.000	ng	0.00
39) Acenaphthene-d10	10.010	164	188182	20.000	ng	0.00
64) Phenanthrene-d10	11.510	188	364742	20.000	ng	0.00
76) Chrysene-d12	14.168	240	242829	20.000	ng	0.00
86) Perylene-d12	15.704	264	178999	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.575	112	661194	144.921	ng	0.02
7) Phenol-d6	6.581	99	816131	140.201	ng	0.00
23) Nitrobenzene-d5	7.522	82	551915	86.683	ng	0.00
42) 2,4,6-Tribromophenol	10.810	330	275202	173.828	ng	0.00
45) 2-Fluorobiphenyl	9.327	172	1050724	81.231	ng	0.00
79) Terphenyl-d14	13.104	244	1374866	92.587	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.851	88	82508	37.933	ng	98
3) Pyridine	3.604	79	224727	37.208	ng	98
4) n-Nitrosodimethylamine	3.575	42	166117	46.639	ng	97
6) Aniline	6.616	93	249059m	32.647	ng	
8) 2-Chlorophenol	6.734	128	255443	49.419	ng	98
9) Benzaldehyde	6.504	77	144188	40.400	ng	97
10) Phenol	6.592	94	308667	47.825	ng	98
11) bis(2-Chloroethyl)ether	6.686	93	240540	44.483	ng	100
12) 1,3-Dichlorobenzene	6.892	146	268069	43.515	ng	98
13) 1,4-Dichlorobenzene	6.969	146	274086	43.787	ng	99
14) 1,2-Dichlorobenzene	7.122	146	266261	46.197	ng	99
15) Benzyl Alcohol	7.098	79	242962	50.903	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.228	45	444599	42.992	ng	100
17) 2-Methylphenol	7.204	107	202498	45.965	ng	99
18) Hexachloroethane	7.469	117	105285	47.019	ng	99
19) n-Nitroso-di-n-propyla...	7.369	70	189662	44.507	ng	97
20) 3+4-Methylphenols	7.357	107	254728m	45.285	ng	
22) Acetophenone	7.369	105	364596	44.628	ng	# 98
24) Nitrobenzene	7.539	77	291541	43.948	ng	98
25) Isophorone	7.781	82	525193	46.234	ng	100
26) 2-Nitrophenol	7.857	139	124809	56.262	ng	98
27) 2,4-Dimethylphenol	7.886	122	226626	51.463	ng	98
28) bis(2-Chloroethoxy)met...	7.986	93	298084	46.201	ng	99
29) 2,4-Dichlorophenol	8.098	162	220147	46.087	ng	96
30) 1,2,4-Trichlorobenzene	8.181	180	250626	42.354	ng	97
31) Naphthalene	8.269	128	717887	41.973	ng	99
32) Benzoic acid	8.016	122	150365	52.714	ng	94
33) 4-Chloroaniline	8.310	127	66241	9.817	ng	97
34) Hexachlorobutadiene	8.380	225	163265	42.609	ng	98
35) Caprolactam	8.692	113	71030m	56.588	ng	
36) 4-Chloro-3-methylphenol	8.792	107	249672	50.310	ng	91
37) 2-Methylnaphthalene	8.963	142	492915	42.534	ng	99
38) 1-Methylnaphthalene	9.063	142	462269	41.136	ng	100
40) 1,2,4,5-Tetrachloroben...	9.127	216	256632	42.024	ng	99
41) Hexachlorocyclopentadiene	9.110	237	338426	123.277	ng	97
43) 2,4,6-Trichlorophenol	9.239	196	172713	49.855	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.275	196	185354	47.581	ng	95
46) 1,1'-Biphenyl	9.427	154	620791	43.229	ng	98
47) 2-Chloronaphthalene	9.457	162	484998	42.101	ng	98
48) 2-Nitroaniline	9.551	65	177913	50.865	ng	98
49) Acenaphthylene	9.874	152	746621	42.516	ng	100
50) Dimethylphthalate	9.733	163	593299	45.251	ng	98
51) 2,6-Dinitrotoluene	9.792	165	131341	50.090	ng	98
52) Acenaphthene	10.045	154	546999m	49.744	ng	
53) 3-Nitroaniline	9.963	138	68425	25.396	ng	92
54) 2,4-Dinitrophenol	10.069	184	143760	105.226	ng	95
55) Dibenzofuran	10.222	168	671038	42.417	ng	99
56) 4-Nitrophenol	10.127	139	216289	115.667	ng	95
57) 2,4-Dinitrotoluene	10.204	165	184060	55.790	ng	88
58) Fluorene	10.563	166	544698	44.117	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.333	232	166987	51.473	ng	99
60) Diethylphthalate	10.433	149	596236	48.161	ng	98
61) 4-Chlorophenyl-phenylether	10.557	204	277058	43.773	ng	96
62) 4-Nitroaniline	10.592	138	130519	48.674	ng	97
63) Azobenzene	10.716	77	583731	43.851	ng	97
65) 4,6-Dinitro-2-methylph...	10.610	198	93866	56.725	ng	84
66) n-Nitrosodiphenylamine	10.674	169	487858	41.976	ng	98
67) 4-Bromophenyl-phenylether	11.051	248	190181	42.542	ng	94
68) Hexachlorobenzene	11.110	284	202610	42.703	ng	98
69) Atrazine	11.204	200	191639	53.554	ng	98
70) Pentachlorophenol	11.304	266	239775	86.547	ng	99
71) Phenanthrene	11.533	178	834900	42.355	ng	100
72) Anthracene	11.586	178	857007	44.239	ng	100
73) Carbazole	11.739	167	764478	46.263	ng	99
74) Di-n-butylphthalate	12.068	149	944967	48.530	ng	99
75) Fluoranthene	12.727	202	925951	44.579	ng	99
77) Benzidine	12.851	184	325361	72.539	ng	97
78) Pyrene	12.963	202	942463	43.990	ng	100
80) Butylbenzylphthalate	13.580	149	349342	48.426	ng	96
81) Benzo(a)anthracene	14.157	228	714641	42.935	ng	99
82) 3,3'-Dichlorobenzidine	14.115	252	109514	24.795	ng	99
83) Chrysene	14.192	228	684758	41.932	ng	99
84) Bis(2-ethylhexyl)phtha...	14.139	149	441324	46.283	ng	# 98
85) Di-n-octyl phthalate	14.768	149	615560	45.882	ng	99
87) Indeno(1,2,3-cd)pyrene	17.280	276	572837	47.648	ng	99
88) Benzo(b)fluoranthene	15.245	252	592913	49.861	ng	99
89) Benzo(k)fluoranthene	15.280	252	550072	45.014	ng	98
90) Benzo(a)pyrene	15.639	252	488310	50.046	ng	98
91) Dibenzo(a,h)anthracene	17.303	278	478539	48.212	ng	97
92) Benzo(g,h,i)perylene	17.762	276	483120	48.768	ng	97

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

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