

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021925\
 Data File : BF141693.D
 Acq On : 20 Feb 2025 00:05
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 02/20/2025
 Supervised By :Jagrut Upadhyay 02/20/2025

Quant Time: Feb 20 02:06:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 16:58:59 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	84080	20.000	ng	-0.01
21) Naphthalene-d8	8.063	136	333578	20.000	ng	-0.01
39) Acenaphthene-d10	9.816	164	175620	20.000	ng	-0.02
64) Phenanthrene-d10	11.298	188	284381	20.000	ng	-0.02
76) Chrysene-d12	13.939	240	159233	20.000	ng	-0.02
86) Perylene-d12	15.380	264	177430	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	468481	87.352	ng	-0.01
7) Phenol-d6	6.422	99	561337	82.811	ng	-0.02
23) Nitrobenzene-d5	7.345	82	511901	80.041	ng	-0.02
42) 2,4,6-Tribromophenol	10.604	330	138958	78.766	ng	-0.02
45) 2-Fluorobiphenyl	9.139	172	964086	84.575	ng	-0.01
79) Terphenyl-d14	12.880	244	840257	89.083	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	2.469	88	94779	43.909	ng	95
3) Pyridine	3.193	79	229986	44.775	ng	97
4) n-Nitrosodimethylamine	3.152	42	129005	37.930	ng	91
6) Aniline	6.445	93	247740	38.961	ng	97
8) 2-Chlorophenol	6.569	128	245940	42.440	ng	98
9) Benzaldehyde	6.334	77	137936	38.293	ng	99
10) Phenol	6.440	94	293175	41.555	ng	96
11) bis(2-Chloroethyl)ether	6.522	93	224785	42.419	ng	99
12) 1,3-Dichlorobenzene	6.722	146	261330	42.715	ng	97
13) 1,4-Dichlorobenzene	6.798	146	266398	42.618	ng	99
14) 1,2-Dichlorobenzene	6.951	146	247985	42.745	ng	100
15) Benzyl Alcohol	6.928	79	207360	39.891	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.057	45	352380	38.846	ng	78
17) 2-Methylphenol	7.045	107	187984	41.452	ng	99
18) Hexachloroethane	7.292	117	98871	42.739	ng	96
19) n-Nitroso-di-n-propyla...	7.198	70	163637	40.358	ng	95
20) 3+4-Methylphenols	7.198	107	239036	41.476	ng	93
22) Acetophenone	7.192	105	337167	40.633	ng	94
24) Nitrobenzene	7.369	77	253453	40.641	ng	98
25) Isophorone	7.604	82	413093	40.509	ng	99
26) 2-Nitrophenol	7.681	139	131772	42.470	ng	94
27) 2,4-Dimethylphenol	7.722	122	149351	41.204	ng	99
28) bis(2-Chloroethoxy)met...	7.816	93	274608	42.025	ng	100
29) 2,4-Dichlorophenol	7.928	162	204723	41.802	ng	99
30) 1,2,4-Trichlorobenzene	8.004	180	221676	41.566	ng	99
31) Naphthalene	8.087	128	718945	41.404	ng	99
32) Benzoic acid	7.851	122	126034	33.475	ng	97
33) 4-Chloroaniline	8.139	127	240828	39.725	ng	98
34) Hexachlorobutadiene	8.198	225	135641	42.366	ng	99
35) Caprolactam	8.510	113	58706	39.370	ng	92
36) 4-Chloro-3-methylphenol	8.628	107	215204	39.669	ng	98
37) 2-Methylnaphthalene	8.775	142	461190	40.816	ng	99
38) 1-Methylnaphthalene	8.875	142	441230	40.318	ng	99
40) 1,2,4,5-Tetrachloroben...	8.939	216	217164	42.741	ng	98
41) Hexachlorocyclopentadiene	8.922	237	49825	29.483	ng	99
43) 2,4,6-Trichlorophenol	9.057	196	136859	41.676	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	147858	41.546	ng	97
46) 1,1'-Biphenyl	9.239	154	576246	42.323	ng	100
47) 2-Chloronaphthalene	9.263	162	423178	42.031	ng	99
48) 2-Nitroaniline	9.363	65	133620	39.545	ng	95
49) Acenaphthylene	9.675	152	624279	41.687	ng	99
50) Dimethylphthalate	9.539	163	501886	42.096	ng	99
51) 2,6-Dinitrotoluene	9.604	165	108613	41.673	ng	98
52) Acenaphthene	9.851	154	407725	40.498	ng	99
53) 3-Nitroaniline	9.775	138	108260	38.593	ng	98
54) 2,4-Dinitrophenol	9.892	184	43173	27.871	ng	91
55) Dibenzofuran	10.022	168	603488	40.838	ng	98
56) 4-Nitrophenol	9.957	139	63683	30.582	ng	92
57) 2,4-Dinitrotoluene	10.010	165	139778	40.286	ng	99
58) Fluorene	10.363	166	467231	40.892	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.145	232	124299	40.567	ng	95
60) Diethylphthalate	10.233	149	481565	41.053	ng	100
61) 4-Chlorophenyl-phenyle...	10.357	204	232006	42.206	ng	96
62) 4-Nitroaniline	10.386	138	104128	36.557	ng	97
63) Azobenzene	10.516	77	465416	39.911	ng	96
65) 4,6-Dinitro-2-methylph...	10.422	198	70281	35.577	ng	92
66) n-Nitrosodiphenylamine	10.475	169	398837	43.812	ng	99
67) 4-Bromophenyl-phenylether	10.845	248	141150	44.410	ng	98
68) Hexachlorobenzene	10.910	284	142480	43.534	ng	98
69) Atrazine	11.004	200	104658	38.322	ng	99
70) Pentachlorophenol	11.110	266	64773	30.952	ng	97
71) Phenanthrene	11.322	178	645237	41.219	ng	100
72) Anthracene	11.375	178	643493	40.893	ng	100
73) Carbazole	11.533	167	549351	38.799	ng	99
74) Di-n-butylphthalate	11.863	149	686027	41.961	ng	100
75) Fluoranthene	12.510	202	602029	36.275	ng	100
77) Benzidine	12.639	184	135198	38.498	ng	99
78) Pyrene	12.739	202	595275	43.915	ng	100
80) Butylbenzylphthalate	13.357	149	219614	46.775	ng	100
81) Benzo(a)anthracene	13.927	228	441236	41.686	ng	99
82) 3,3'-Dichlorobenzidine	13.892	252	137095	45.215	ng	100
83) Chrysene	13.963	228	415159	42.478	ng	100
84) Bis(2-ethylhexyl)phtha...	13.916	149	286347	54.076	ng	100
85) Di-n-octyl phthalate	14.527	149	432533	57.257	ng	97
87) Indeno(1,2,3-cd)pyrene	16.815	276	514199	46.902	ng	99
88) Benzo(b)fluoranthene	14.963	252	431312	36.203	ng	100
89) Benzo(k)fluoranthene	14.992	252	419957m	41.281	ng	
90) Benzo(a)pyrene	15.321	252	394420	40.915	ng	99
91) Dibenzo(a,h)anthracene	16.827	278	419567	47.231	ng	99
92) Benzo(g,h,i)perylene	17.245	276	429072	46.156	ng	100

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

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