

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121222\
 Data File : BF131594.D
 Acq On : 12 Dec 2022 15:59
 Operator : CG\JU
 Sample : N5964-09MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-518-COMP-05MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/13/2022
 Supervised By : Jagrut Upadhyay 12/13/2022

Quant Time: Dec 13 04:15:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 13 04:12:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	62043	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	223107	20.000	ng	# 0.00
39) Acenaphthene-d10	9.951	164	128415	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	242568	20.000	ng	0.00
76) Chrysene-d12	14.104	240	137964	20.000	ng	# 0.00
86) Perylene-d12	15.615	264	136691	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	438171	117.821	ng	0.01
7) Phenol-d6	6.534	99	571053	117.177	ng	0.00
23) Nitrobenzene-d5	7.463	82	390625	66.194	ng	0.00
42) 2,4,6-Tribromophenol	10.751	330	168545	118.134	ng	0.00
45) 2-Fluorobiphenyl	9.269	172	669263	65.713	ng	0.00
79) Terphenyl-d14	13.039	244	671790	66.778	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.734	88	73584	43.835	ng	98
3) Pyridine	3.487	79	196915	42.251	ng	98
4) n-Nitrosodimethylamine	3.457	42	109878	42.749	ng	100
6) Aniline	6.557	93	231475	39.715	ng	99
8) 2-Chlorophenol	6.681	128	185122	46.972	ng	93
9) Benzaldehyde	6.445	77	102457	30.689	ng	98
10) Phenol	6.545	94	265934m	48.921	ng	
11) bis(2-Chloroethyl)ether	6.628	93	189367	46.472	ng	98
12) 1,3-Dichlorobenzene	6.834	146	207119	45.328	ng	97
13) 1,4-Dichlorobenzene	6.910	146	212678	46.144	ng	98
14) 1,2-Dichlorobenzene	7.063	146	201939	45.866	ng	97
15) Benzyl Alcohol	7.039	79	203734	45.600	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.169	45	228032	43.322	ng	97
17) 2-Methylphenol	7.151	107	161110	45.359	ng	98
18) Hexachloroethane	7.404	117	84986	45.710	ng	91
19) n-Nitroso-di-n-propyla...	7.310	70	160860	44.019	ng	99
20) 3+4-Methylphenols	7.304	107	220379	45.553	ng	# 82
22) Acetophenone	7.304	105	308462	44.119	ng	94
24) Nitrobenzene	7.481	77	250623	42.108	ng	95
25) Isophorone	7.722	82	437879	46.066	ng	99
26) 2-Nitrophenol	7.798	139	98972	48.762	ng	93
27) 2,4-Dimethylphenol	7.834	122	167339	53.789	ng	99
28) bis(2-Chloroethoxy)met...	7.934	93	241452	49.103	ng	100
29) 2,4-Dichlorophenol	8.039	162	171182	44.501	ng	96
30) 1,2,4-Trichlorobenzene	8.122	180	202080	42.837	ng	98
31) Naphthalene	8.204	128	531422	42.290	ng	100
32) Benzoic acid	7.969	122	110208	48.915	ng	97
33) 4-Chloroaniline	8.257	127	115830	24.591	ng	96
34) Hexachlorobutadiene	8.316	225	135592	40.784	ng	99
35) Caprolactam	8.639	113	48425m	48.665	ng	
36) 4-Chloro-3-methylphenol	8.745	107	194783	45.094	ng	98
37) 2-Methylnaphthalene	8.898	142	367643	42.780	ng	99
38) 1-Methylnaphthalene	8.998	142	346365	41.985	ng	98
40) 1,2,4,5-Tetrachloroben...	9.069	216	217083	45.667	ng	98
41) Hexachlorocyclopentadiene	9.051	237	266625	110.503	ng	98
43) 2,4,6-Trichlorophenol	9.180	196	135638	46.131	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.228	196	141599	44.780	ng	99
46) 1,1'-Biphenyl	9.369	154	471619	46.450	ng	99
47) 2-Chloronaphthalene	9.392	162	375381	45.378	ng	98
48) 2-Nitroaniline	9.498	65	129966	44.352	ng	89
49) Acenaphthylene	9.816	152	558245	45.961	ng	99
50) Dimethylphthalate	9.675	163	436377	44.577	ng	99
51) 2,6-Dinitrotoluene	9.739	165	96751	47.874	ng	# 82
52) Acenaphthene	9.986	154	407052m	50.026	ng	
53) 3-Nitroaniline	9.910	138	57850	29.835	ng	99
54) 2,4-Dinitrophenol	10.016	184	101693	87.497	ng	92
55) Dibenzofuran	10.163	168	520865	45.305	ng	96
56) 4-Nitrophenol	10.080	139	136679	101.467	ng	89
57) 2,4-Dinitrotoluene	10.145	165	129921	48.462	ng	96
58) Fluorene	10.504	166	412591	44.144	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.280	232	118889m	44.400	ng	
60) Diethylphthalate	10.380	149	415823	44.264	ng	99
61) 4-Chlorophenyl-phenyle...	10.498	204	227592	44.064	ng	93
62) 4-Nitroaniline	10.533	138	86801	45.174	ng	95
63) Azobenzene	10.657	77	455453	42.508	ng	98
65) 4,6-Dinitro-2-methylph...	10.557	198	68512	44.594	ng	78
66) n-Nitrosodiphenylamine	10.616	169	346797	44.223	ng	98
67) 4-Bromophenyl-phenylether	10.986	248	134628	43.144	ng	95
68) Hexachlorobenzene	11.051	284	146839	43.504	ng	# 93
69) Atrazine	11.151	200	130255	52.305	ng	96
70) Pentachlorophenol	11.251	266	153533	88.358	ng	97
71) Phenanthrene	11.474	178	583967	43.191	ng	100
72) Anthracene	11.527	178	599256	45.556	ng	99
73) Carbazole	11.680	167	492146	45.836	ng	99
74) Di-n-butylphthalate	12.010	149	601017	45.627	ng	100
75) Fluoranthene	12.669	202	606686	43.576	ng	99
77) Benzidine	12.792	184	227648	61.583	ng	98
78) Pyrene	12.898	202	597194	43.898	ng	99
80) Butylbenzylphthalate	13.515	149	209317	53.168	ng	99
81) Benzo(a)anthracene	14.092	228	452251	45.392	ng	99
82) 3,3'-Dichlorobenzidine	14.057	252	101401	35.772	ng	97
83) Chrysene	14.133	228	424566	44.823	ng	98
84) Bis(2-ethylhexyl)phtha...	14.074	149	267462	49.750	ng	99
85) Di-n-octyl phthalate	14.698	149	396487	45.429	ng	98
87) Indeno(1,2,3-cd)pyrene	17.156	276	451956	40.514	ng	99
88) Benzo(b)fluoranthene	15.168	252	406024	45.068	ng	100
89) Benzo(k)fluoranthene	15.198	252	412501	43.868	ng	99
90) Benzo(a)pyrene	15.551	252	356651	46.388	ng	100
91) Dibenzo(a,h)anthracene	17.180	278	384843	41.917	ng	98
92) Benzo(g,h,i)perylene	17.627	276	380780	41.267	ng	99

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

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