

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032322\
 Data File : BF127485.D
 Acq On : 23 Mar 2022 17:25
 Operator : CG\JU
 Sample : N2058-02MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MHAMSD

Quant Time: Mar 24 02:12:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 22 02:16:33 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/24/2022
 Supervised By : Jagrut Upadhyay 03/24/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	88282	20.000	ng	0.00
21) Naphthalene-d8	8.134	136	329371	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	190693	20.000	ng	0.00
64) Phenanthrene-d10	11.380	188	304968	20.000	ng	0.00
76) Chrysene-d12	14.027	240	183594	20.000	ng	0.00
86) Perylene-d12	15.510	264	171721	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	747967	135.239	ng	0.02
7) Phenol-d6	6.493	99	962530	127.318	ng	0.00
23) Nitrobenzene-d5	7.422	82	743003	96.427	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	265227	130.566	ng	0.00
45) 2-Fluorobiphenyl	9.210	172	1211383	92.762	ng	0.00
79) Terphenyl-d14	12.963	244	1062225	94.714	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.834	88	124406	51.214	ng	# 79
3) Pyridine	3.552	79	221196	32.263	ng	96
4) n-Nitrosodimethylamine	3.499	42	204866	54.423	ng	98
6) Aniline	6.522	93	85560	9.484	ng	# 54
8) 2-Chlorophenol	6.640	128	312591	55.308	ng	97
9) Benzaldehyde	6.410	77	144641	36.812	ng	99
10) Phenol	6.504	94	381112	46.290	ng	99
11) bis(2-Chloroethyl)ether	6.587	93	348131	57.787	ng	98
12) 1,3-Dichlorobenzene	6.787	146	326640	51.384	ng	99
13) 1,4-Dichlorobenzene	6.863	146	328743	51.412	ng	97
14) 1,2-Dichlorobenzene	7.016	146	315554	50.668	ng	99
15) Benzyl Alcohol	6.998	79	254031	46.236	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.122	45	576677	53.196	ng	82
17) 2-Methylphenol	7.110	107	266042	55.776	ng	94
18) Hexachloroethane	7.351	117	126280	53.944	ng	93
19) n-Nitroso-di-n-propyla...	7.269	70	248035	54.734	ng	98
20) 3+4-Methylphenols	7.263	107	314577	53.471	ng	# 83
22) Acetophenone	7.263	105	434539	51.611	ng	96
24) Nitrobenzene	7.440	77	391223	53.239	ng	98
25) Isophorone	7.675	82	704555	57.410	ng	99
26) 2-Nitrophenol	7.751	139	162301	52.662	ng	98
27) 2,4-Dimethylphenol	7.787	122	284595	61.389	ng	96
28) bis(2-Chloroethoxy)met...	7.881	93	414009	51.932	ng	99
29) 2,4-Dichlorophenol	7.998	162	275785	52.402	ng	97
30) 1,2,4-Trichlorobenzene	8.069	180	303418	49.475	ng	100
31) Naphthalene	8.151	128	903157	51.648	ng	99
32) Benzoic acid	7.951	122	147367	54.390	ng	92
33) 4-Chloroaniline	8.222	127	30739	4.536	ng	98
34) Hexachlorobutadiene	8.257	225	190616	48.947	ng	99
35) Caprolactam	8.598	113	69618m	42.908	ng	
36) 4-Chloro-3-methylphenol	8.698	107	296245	51.881	ng	97
37) 2-Methylnaphthalene	8.845	142	592138	52.662	ng	100
38) 1-Methylnaphthalene	8.945	142	562392	50.582	ng	99
40) 1,2,4,5-Tetrachloroben...	9.010	216	304008	50.140	ng	99
41) Hexachlorocyclopentadiene	8.992	237	245149	102.529	ng	99
43) 2,4,6-Trichlorophenol	9.128	196	188121	51.329	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.181	196	200242	53.076	ng	94
46) 1,1'-Biphenyl	9.310	154	727249	50.495	ng	99
47) 2-Chloronaphthalene	9.339	162	568514	50.127	ng	99
48) 2-Nitroaniline	9.445	65	203891	46.907	ng	96
49) Acenaphthylene	9.751	152	897938	52.371	ng	99
50) Dimethylphthalate	9.616	163	689006	51.937	ng	100
51) 2,6-Dinitrotoluene	9.686	165	145561	52.271	ng	# 83
52) Acenaphthene	9.922	154	510175	50.998	ng	99
53) 3-Nitroaniline	9.857	138	21262	6.960	ng	97
54) 2,4-Dinitrophenol	9.969	184	143035	99.404	ng	# 89
55) Dibenzofuran	10.098	168	740446	51.797	ng	99
56) 4-Nitrophenol	10.034	139	118686	76.338	ng	96
57) 2,4-Dinitrotoluene	10.092	165	173838	48.499	ng	97
58) Fluorene	10.439	166	596356	51.828	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.222	232	165216	49.531	ng	97
60) Diethylphthalate	10.310	149	585966	48.360	ng	99
61) 4-Chlorophenyl-phenyle...	10.428	204	320255	46.709	ng	99
62) 4-Nitroaniline	10.475	138	85529	29.892	ng	97
63) Azobenzene	10.592	77	679157	47.503	ng	98
65) 4,6-Dinitro-2-methylph...	10.504	198	95676	50.985	ng	89
66) n-Nitrosodiphenylamine	10.551	169	401653	43.266	ng	99
67) 4-Bromophenyl-phenylether	10.916	248	192725	54.128	ng	93
68) Hexachlorobenzene	10.986	284	213542	54.944	ng	97
69) Atrazine	11.075	200	93379	30.215	ng	96
70) Pentachlorophenol	11.186	266	171597	111.458	ng	98
71) Phenanthrene	11.404	178	833013	52.393	ng	99
72) Anthracene	11.457	178	822252	52.114	ng	99
73) Carbazole	11.616	167	561512	37.214	ng	99
74) Di-n-butylphthalate	11.933	149	814992	52.181	ng	100
75) Fluoranthene	12.598	202	784249	44.846	ng	99
77) Benzidine	12.745	184	8771	2.004	ng	# 91
78) Pyrene	12.827	202	772828	49.048	ng	99
80) Butylbenzylphthalate	13.433	149	260737	50.279	ng	94
81) Benzo(a)anthracene	14.016	228	635159	51.732	ng	98
82) 3,3'-Dichlorobenzidine	13.980	252	15860	4.395	ng	97
83) Chrysene	14.051	228	605328	52.065	ng	100
84) Bis(2-ethylhexyl)phtha...	13.980	149	343909	57.370	ng	98
85) Di-n-octyl phthalate	14.598	149	591613	74.164	ng	99
87) Indeno(1,2,3-cd)pyrene	17.010	276	584251	50.281	ng	99
88) Benzo(b)fluoranthene	15.074	252	624729	57.699	ng	98
89) Benzo(k)fluoranthene	15.104	252	625242	55.704	ng	99
90) Benzo(a)pyrene	15.445	252	604974	68.083	ng	98
91) Dibenzo(a,h)anthracene	17.021	278	533973	55.661	ng	99
92) Benzo(g,h,i)perylene	17.462	276	485299	48.942	ng	99

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

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