

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102124\
 Data File : BF139914.D
 Acq On : 21 Oct 2024 20:32
 Operator : RC/JU
 Sample : P4419-01MSD
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IB-1MSD

Quant Time: Oct 22 01:20:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/22/2024
 Supervised By :mohammad ahmed 10/23/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	130734	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	435954	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	207044	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	353526	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	219663	20.000	ng	0.00
86) Perylene-d12	15.539	264	191968	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	864546	103.526	ng	0.02
7) Phenol-d6	6.510	99	1068913	98.828	ng	0.00
23) Nitrobenzene-d5	7.451	82	637237	81.013	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	223502	115.408	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	899868	71.831	ng	0.00
79) Terphenyl-d14	13.004	244	826375	61.301	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.746	88	188299	48.361	ng	97
3) Pyridine	3.499	79	477302	49.455	ng	100
4) n-Nitrosodimethylamine	3.452	42	273365	52.719	ng	100
6) Aniline	6.551	93	361812	35.893	ng	99
8) 2-Chlorophenol	6.675	128	442375	51.817	ng	99
9) Benzaldehyde	6.440	77	110558	18.170	ng	98
10) Phenol	6.528	94	581563m	52.155	ng	
11) bis(2-Chloroethyl)ether	6.628	93	453156	52.579	ng	100
12) 1,3-Dichlorobenzene	6.834	146	479386	48.917	ng	99
13) 1,4-Dichlorobenzene	6.910	146	482780	49.309	ng	99
14) 1,2-Dichlorobenzene	7.063	146	456437	49.950	ng	99
15) Benzyl Alcohol	7.028	79	405071	51.056	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.163	45	731039	49.744	ng	99
17) 2-Methylphenol	7.140	107	357583	49.120	ng	100
18) Hexachloroethane	7.404	117	147194	42.159	ng	98
19) n-Nitroso-di-n-propyla...	7.304	70	319944	48.773	ng	97
20) 3+4-Methylphenols	7.292	107	445899	47.888	ng	# 84
22) Acetophenone	7.298	105	589724	55.228	ng	99
24) Nitrobenzene	7.475	77	462958	53.682	ng	99
25) Isophorone	7.710	82	822101	55.066	ng	99
26) 2-Nitrophenol	7.787	139	189017	58.097	ng	98
27) 2,4-Dimethylphenol	7.822	122	340338	62.735	ng	100
28) bis(2-Chloroethoxy)met...	7.922	93	491123	54.296	ng	98
29) 2,4-Dichlorophenol	8.028	162	325630	52.698	ng	99
30) 1,2,4-Trichlorobenzene	8.116	180	345488	50.529	ng	99
31) Naphthalene	8.198	128	1152723	51.269	ng	100
32) Benzoic acid	7.922	122	234230	49.503	ng	96
33) 4-Chloroaniline	8.239	127	169402	22.096	ng	99
34) Hexachlorobutadiene	8.316	225	226524	52.729	ng	100
35) Caprolactam	8.598	113	116327	59.206	ng	# 84
36) 4-Chloro-3-methylphenol	8.716	107	332636	48.616	ng	99
37) 2-Methylnaphthalene	8.886	142	679816	49.369	ng	100
38) 1-Methylnaphthalene	8.986	142	628892	46.550	ng	99
40) 1,2,4,5-Tetrachloroben...	9.051	216	317228	56.793	ng	99
41) Hexachlorocyclopentadiene	9.039	237	105930	54.503	ng	98
43) 2,4,6-Trichlorophenol	9.163	196	219318	55.458	ng	99

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44) 2,4,5-Trichlorophenol	9.204	196	213935	52.434	ng	96
46) 1,1'-Biphenyl	9.351	154	780261	54.135	ng	100
47) 2-Chloronaphthalene	9.375	162	604418	52.066	ng	99
48) 2-Nitroaniline	9.469	65	214962	60.545	ng	94
49) Acenaphthylene	9.792	152	913331	54.421	ng	99
50) Dimethylphthalate	9.651	163	699542	54.244	ng	100
51) 2,6-Dinitrotoluene	9.710	165	137647	49.295	ng	98
52) Acenaphthene	9.969	154	565907	52.126	ng	99
53) 3-Nitroaniline	9.875	138	142625	49.530	ng	100
54) 2,4-Dinitrophenol	9.981	184	35986	35.015	ng	81
55) Dibenzofuran	10.139	168	806946	51.805	ng	99
56) 4-Nitrophenol	10.028	139	258658	116.755	ng	97
57) 2,4-Dinitrotoluene	10.110	165	175165	51.012	ng	# 99
58) Fluorene	10.481	166	620230	52.278	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.251	232	171768	53.702	ng	# 97
60) Diethylphthalate	10.351	149	670868	52.981	ng	99
61) 4-Chlorophenyl-phenyle...	10.475	204	303779	50.703	ng	100
62) 4-Nitroaniline	10.486	138	143901	52.912	ng	99
63) Azobenzene	10.633	77	716946	53.408	ng	93
65) 4,6-Dinitro-2-methylph...	10.522	198	27186	18.853	ng	# 64
66) n-Nitrosodiphenylamine	10.586	169	614128	58.041	ng	97
67) 4-Bromophenyl-phenylether	10.963	248	194148	53.308	ng	99
68) Hexachlorobenzene	11.028	284	211775	51.703	ng	99
69) Atrazine	11.116	200	186610	65.107	ng	99
70) Pentachlorophenol	11.222	266	276377	111.178	ng	99
71) Phenanthrene	11.445	178	974332	58.346	ng	99
72) Anthracene	11.498	178	999042	61.317	ng	99
73) Carbazole	11.651	167	850430	56.123	ng	99
74) Di-n-butylphthalate	11.980	149	1026526	58.636	ng	100
75) Fluoranthene	12.633	202	1064175	63.038	ng	99
77) Benzidine	12.751	184	432650	119.781	ng	99
78) Pyrene	12.863	202	1081259	56.234	ng	100
80) Butylbenzylphthalate	13.474	149	418659	72.627	ng	97
81) Benzo(a)anthracene	14.051	228	842341	58.908	ng	99
82) 3,3'-Dichlorobenzidine	14.010	252	247304	59.317	ng	100
83) Chrysene	14.086	228	719453	54.875	ng	99
84) Bis(2-ethylhexyl)phtha...	14.039	149	526835	81.459	ng	# 98
85) Di-n-octyl phthalate	14.651	149	802185	68.192	ng	98
87) Indeno(1,2,3-cd)pyrene	17.033	276	582408	47.159	ng	98
88) Benzo(b)fluoranthene	15.104	252	641069	54.821	ng	99
89) Benzo(k)fluoranthene	15.133	252	622871	61.762	ng	100
90) Benzo(a)pyrene	15.474	252	589245	61.239	ng	99
91) Dibenzo(a,h)anthracene	17.057	278	474682	46.075	ng	99
92) Benzo(g,h,i)perylene	17.486	276	418831	40.699	ng	98

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

