

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010423\
 Data File : BF131901.D
 Acq On : 04 Jan 2023 16:17
 Operator : CG\JU
 Sample : 01005-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-10323MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/05/2023
 Supervised By : Jagrut Upadhyay 01/05/2023

Quant Time: Jan 05 05:42:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 30 02:06:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	71344	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	253565	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	135856	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	221819	20.000	ng	0.00
76) Chrysene-d12	14.016	240	134129	20.000	ng	0.00
86) Perylene-d12	15.492	264	160208	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	498518	115.799	ng	0.01
7) Phenol-d6	6.457	99	655154	115.833	ng	0.00
23) Nitrobenzene-d5	7.387	82	452493	71.248	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	160052	108.430	ng	0.00
45) 2-Fluorobiphenyl	9.187	172	727876	73.386	ng	0.00
79) Terphenyl-d14	12.951	244	510821	64.420	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.581	88	81671	40.910	ng	97
3) Pyridine	3.310	79	227855	40.599	ng	98
4) n-Nitrosodimethylamine	3.287	42	116538	44.162	ng	99
6) Aniline	6.481	93	228099	33.253	ng	99
8) 2-Chlorophenol	6.598	128	206947	45.627	ng	94
9) Benzaldehyde	6.363	77	135674	37.384	ng	99
10) Phenol	6.469	94	292958	43.494	ng	92
11) bis(2-Chloroethyl)ether	6.551	93	221034	45.890	ng	99
12) 1,3-Dichlorobenzene	6.751	146	231295	44.624	ng	98
13) 1,4-Dichlorobenzene	6.834	146	235545	45.093	ng	96
14) 1,2-Dichlorobenzene	6.987	146	223842	45.605	ng	99
15) Benzyl Alcohol	6.963	79	226614	45.689	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.093	45	278831	45.363	ng	96
17) 2-Methylphenol	7.081	107	185437	44.722	ng	97
18) Hexachloroethane	7.328	117	93358	44.260	ng	92
19) n-Nitroso-di-n-propyla...	7.234	70	177854	44.524	ng	98
20) 3+4-Methylphenols	7.234	107	234161	43.281	ng	99
22) Acetophenone	7.234	105	340056	45.745	ng	# 97
24) Nitrobenzene	7.404	77	278820	43.807	ng	97
25) Isophorone	7.645	82	488237	47.074	ng	99
26) 2-Nitrophenol	7.722	139	110056	44.473	ng	93
27) 2,4-Dimethylphenol	7.763	122	187095	53.003	ng	97
28) bis(2-Chloroethoxy)met...	7.857	93	272057	48.895	ng	100
29) 2,4-Dichlorophenol	7.969	162	185923	43.718	ng	99
30) 1,2,4-Trichlorobenzene	8.045	180	219623	43.556	ng	99
31) Naphthalene	8.128	128	592059	42.635	ng	99
32) Benzoic acid	7.904	122	112340	43.542	ng	91
33) 4-Chloroaniline	8.181	127	86753	16.282	ng	98
34) Hexachlorobutadiene	8.240	225	147380	42.723	ng	99
35) Caprolactam	8.563	113	53247m	42.288	ng	
36) 4-Chloro-3-methylphenol	8.675	107	209718	43.632	ng	98
37) 2-Methylnaphthalene	8.822	142	395947	42.484	ng	99
38) 1-Methylnaphthalene	8.922	142	362787	40.075	ng	99
40) 1,2,4,5-Tetrachloroben...	8.987	216	228175	47.833	ng	99
41) Hexachlorocyclopentadiene	8.969	237	182485	85.074	ng	97
43) 2,4,6-Trichlorophenol	9.104	196	139212	45.566	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010423\
 Data File : BF131901.D
 Acq On : 04 Jan 2023 16:17
 Operator : CG\JU
 Sample : 01005-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-10323MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/05/2023
 Supervised By : Jagrut Upadhyay 01/05/2023

Quant Time: Jan 05 05:42:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 30 02:06:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.145	196	145284	43.991	ng	99
46) 1,1'-Biphenyl	9.287	154	496959	47.779	ng	100
47) 2-Chloronaphthalene	9.316	162	392571	45.621	ng	100
48) 2-Nitroaniline	9.416	65	135950	43.103	ng	97
49) Acenaphthylene	9.734	152	580489	45.304	ng	99
50) Dimethylphthalate	9.592	163	471350	45.860	ng	99
51) 2,6-Dinitrotoluene	9.663	165	99628	44.778	ng	90
52) Acenaphthene	9.904	154	336777	43.550	ng	100
53) 3-Nitroaniline	9.828	138	60299	26.598	ng	97
54) 2,4-Dinitrophenol	9.939	184	95535	74.614	ng	98
55) Dibenzofuran	10.075	168	524234	44.000	ng	99
56) 4-Nitrophenol	10.004	139	121212	76.325	ng	93
57) 2,4-Dinitrotoluene	10.069	165	130119	42.763	ng	96
58) Fluorene	10.422	166	403420	41.476	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.198	232	112784	41.615	ng	# 88
60) Diethylphthalate	10.292	149	440010	44.481	ng	98
61) 4-Chlorophenyl-phenyle...	10.410	204	224587	44.417	ng	99
62) 4-Nitroaniline	10.451	138	83065	35.157	ng	96
63) Azobenzene	10.575	77	466686	42.509	ng	99
65) 4,6-Dinitro-2-methylph...	10.475	198	62781	41.569	ng	86
66) n-Nitrosodiphenylamine	10.534	169	343592	50.779	ng	98
67) 4-Bromophenyl-phenylether	10.904	248	126964	49.385	ng	97
68) Hexachlorobenzene	10.969	284	133480	47.263	ng	# 89
69) Atrazine	11.063	200	120316	54.011	ng	97
70) Pentachlorophenol	11.169	266	120270	86.063	ng	98
71) Phenanthrene	11.386	178	561957	46.069	ng	99
72) Anthracene	11.439	178	534077	44.757	ng	99
73) Carbazole	11.598	167	438079	41.586	ng	99
74) Di-n-butylphthalate	11.922	149	578576	45.439	ng	99
75) Fluoranthene	12.580	202	553985	38.990	ng	100
77) Benzidine	12.704	184	182222	82.669	ng	99
78) Pyrene	12.810	202	543963	48.386	ng	99
80) Butylbenzylphthalate	13.427	149	187811	45.040	ng	93
81) Benzo(a)anthracene	14.004	228	466503	48.038	ng	99
82) 3,3'-Dichlorobenzidine	13.969	252	115011	40.975	ng	97
83) Chrysene	14.045	228	440906	48.315	ng	98
84) Bis(2-ethylhexyl)phtha...	13.986	149	249744	49.190	ng	# 97
85) Di-n-octyl phthalate	14.604	149	449727	65.800	ng	99
87) Indeno(1,2,3-cd)pyrene	16.980	276	508989	50.008	ng	99
88) Benzo(b)fluoranthene	15.063	252	497913	42.386	ng	100
89) Benzo(k)fluoranthene	15.092	252	469232	41.341	ng	98
90) Benzo(a)pyrene	15.433	252	438615	48.228	ng	97
91) Dibenzo(a,h)anthracene	16.998	278	425738	51.065	ng	98
92) Benzo(g,h,i)perylene	17.433	276	391397	40.600	ng	96

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

