

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020824\
 Data File : BF137337.D
 Acq On : 08 Feb 2024 14:57
 Operator : CG\JU
 Sample : PB158865BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB158865BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/09/2024
 Supervised By :mohammad ahmed 02/09/2024

Quant Time: Feb 08 15:33:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 04:56:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	102217	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	400856	20.000	ng	#-0.01
39) Acenaphthene-d10	9.822	164	223152	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	408780	20.000	ng	0.00
76) Chrysene-d12	13.968	240	241393	20.000	ng	0.00
86) Perylene-d12	15.445	264	146725	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	813208	131.276	ng	0.01
7) Phenol-d6	6.428	99	1066035	120.587	ng	0.00
23) Nitrobenzene-d5	7.351	82	716271	80.000	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	293664	116.525	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	1226964	75.584	ng	0.00
79) Terphenyl-d14	12.910	244	1331985	75.013	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.646	88	115887	34.963	ng	97
3) Pyridine	3.387	79	274155	41.305	ng	98
4) n-Nitrosodimethylamine	3.357	42	145069	43.584	ng	90
6) Aniline	6.451	93	321765	34.001	ng	98
8) 2-Chlorophenol	6.569	128	306120	44.196	ng	99
9) Benzaldehyde	6.340	77	89657	26.631	ng	96
10) Phenol	6.440	94	414856	42.041	ng	97
11) bis(2-Chloroethyl)ether	6.528	93	293407	44.492	ng	99
12) 1,3-Dichlorobenzene	6.728	146	332609	40.957	ng	96
13) 1,4-Dichlorobenzene	6.804	146	345328	41.117	ng	96
14) 1,2-Dichlorobenzene	6.951	146	324006	41.261	ng	99
15) Benzyl Alcohol	6.934	79	290433	46.243	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.057	45	512887	43.175	ng	98
17) 2-Methylphenol	7.039	107	261453	45.183	ng	97
18) Hexachloroethane	7.292	117	123336	40.717	ng	95
19) n-Nitroso-di-n-propyla...	7.204	70	242517	40.010	ng	99
20) 3+4-Methylphenols	7.192	107	308016	42.048	ng	94
22) Acetophenone	7.198	105	431616	39.995	ng	# 97
24) Nitrobenzene	7.369	77	364037	41.891	ng	96
25) Isophorone	7.604	82	678237	41.040	ng	97
26) 2-Nitrophenol	7.681	139	153390	45.716	ng	93
27) 2,4-Dimethylphenol	7.722	122	212177	44.918	ng	97
28) bis(2-Chloroethoxy)met...	7.822	93	382652	43.602	ng	99
29) 2,4-Dichlorophenol	7.922	162	258205	44.063	ng	99
30) 1,2,4-Trichlorobenzene	8.004	180	304240	41.798	ng	98
31) Naphthalene	8.086	128	888319	40.992	ng	99
32) Benzoic acid	7.851	122	162872	48.779	ng	93
33) 4-Chloroaniline	8.139	127	110985	14.889	ng	95
34) Hexachlorobutadiene	8.204	225	197148	40.734	ng	94
35) Caprolactam	8.510	113	80642m	44.258	ng	
36) 4-Chloro-3-methylphenol	8.622	107	286034	41.327	ng	95
37) 2-Methylnaphthalene	8.775	142	586862	40.394	ng	100
38) 1-Methylnaphthalene	8.875	142	563451	41.755	ng	100
40) 1,2,4,5-Tetrachloroben...	8.945	216	310844	41.630	ng	99
41) Hexachlorocyclopentadiene	8.928	237	380352	106.950	ng	98
43) 2,4,6-Trichlorophenol	9.057	196	189723	44.797	ng	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020824\
 Data File : BF137337.D
 Acq On : 08 Feb 2024 14:57
 Operator : CG\JU
 Sample : PB158865BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB158865BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/09/2024
 Supervised By :mohammad ahmed 02/09/2024

Quant Time: Feb 08 15:33:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 04:56:22 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.098	196	201872	43.607	ng	96
46) 1,1'-Biphenyl	9.245	154	734319	41.359	ng	99
47) 2-Chloronaphthalene	9.269	162	542038	41.253	ng	99
48) 2-Nitroaniline	9.369	65	196791	46.274	ng	96
49) Acenaphthylene	9.686	152	861572	40.959	ng	99
50) Dimethylphthalate	9.551	163	628708	41.704	ng	100
51) 2,6-Dinitrotoluene	9.610	165	144555	44.967	ng	92
52) Acenaphthene	9.857	154	513631	41.547	ng	99
53) 3-Nitroaniline	9.780	138	77393	26.752	ng	94
54) 2,4-Dinitrophenol	9.886	184	139682	95.062	ng	85
55) Dibenzofuran	10.028	168	787117	40.019	ng	98
56) 4-Nitrophenol	9.945	139	201928	97.238	ng	95
57) 2,4-Dinitrotoluene	10.016	165	184999	44.352	ng	89
58) Fluorene	10.375	166	594829	38.577	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.145	232	185768m	43.128	ng	
60) Diethylphthalate	10.251	149	601776	40.540	ng	98
61) 4-Chlorophenyl-phenyle...	10.369	204	296535	38.690	ng	96
62) 4-Nitroaniline	10.404	138	122316m	44.172	ng	
63) Azobenzene	10.527	77	658269	40.042	ng	96
65) 4,6-Dinitro-2-methylph...	10.422	198	104884	48.387	ng	89
66) n-Nitrosodiphenylamine	10.486	169	532371	42.832	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	194796	43.338	ng	96
68) Hexachlorobenzene	10.916	284	210777	41.982	ng	99
69) Atrazine	11.022	200	183145	44.635	ng	97
70) Pentachlorophenol	11.116	266	265969	92.413	ng	99
71) Phenanthrene	11.339	178	885768	41.391	ng	100
72) Anthracene	11.392	178	919593	41.228	ng	99
73) Carbazole	11.551	167	797231	42.767	ng	99
74) Di-n-butylphthalate	11.886	149	874086	43.378	ng	99
75) Fluoranthene	12.533	202	904785	40.404	ng	99
77) Benzidine	12.663	184	223087	51.611	ng	99
78) Pyrene	12.763	202	929105	40.051	ng	99
80) Butylbenzylphthalate	13.392	149	312251	52.618	ng	94
81) Benzo(a)anthracene	13.957	228	702053	41.066	ng	99
82) 3,3'-Dichlorobenzidine	13.927	252	171672	36.858	ng	98
83) Chrysene	13.992	228	721007	42.899	ng	99
84) Bis(2-ethylhexyl)phtha...	13.963	149	386420	55.970	ng	# 97
85) Di-n-octyl phthalate	14.580	149	633530	57.711	ng	100
87) Indeno(1,2,3-cd)pyrene	16.945	276	472697	45.856	ng	99
88) Benzo(b)fluoranthene	15.015	252	616107	68.279	ng	98
89) Benzo(k)fluoranthene	15.045	252	619343	64.881	ng	100
90) Benzo(a)pyrene	15.380	252	436673	51.424	ng	99
91) Dibenzo(a,h)anthracene	16.974	278	384258	46.522	ng	98
92) Benzo(g,h,i)perylene	17.404	276	374414	46.186	ng	99

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

