

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120123\
 Data File : BF136332.D
 Acq On : 01 Dec 2023 12:08
 Operator : CG\JU
 Sample : PB157138BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB157138BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/04/2023
 Supervised By :mohammad ahmed 12/05/2023

Quant Time: Dec 01 13:16:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 03:04:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	138172	20.000	ng	0.00
21) Naphthalene-d8	8.087	136	613120	20.000	ng	0.00
39) Acenaphthene-d10	9.845	164	314144	20.000	ng	0.00
64) Phenanthrene-d10	11.339	188	537508	20.000	ng	0.00
76) Chrysene-d12	13.992	240	239799	20.000	ng	0.00
86) Perylene-d12	15.474	264	260106	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	1123372	113.161	ng	0.02
7) Phenol-d6	6.451	99	1396891	112.537	ng	0.00
23) Nitrobenzene-d5	7.375	82	834223	71.900	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	429646	111.867	ng	0.00
45) 2-Fluorobiphenyl	9.169	172	1757454	71.458	ng	0.00
79) Terphenyl-d14	12.933	244	1515189	72.884	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.681	88	117902	31.389	ng	95
3) Pyridine	3.440	79	286544	30.903	ng	95
4) n-Nitrosodimethylamine	3.405	42	165614	35.335	ng	92
6) Aniline	6.475	93	370899	31.614	ng	97
8) 2-Chlorophenol	6.598	128	443858	41.063	ng	91
9) Benzaldehyde	6.357	77	21281	3.076	ng	95
10) Phenol	6.463	94	521735	39.179	ng	98
11) bis(2-Chloroethyl)ether	6.551	93	392130	40.139	ng	94
12) 1,3-Dichlorobenzene	6.751	146	457547	37.300	ng	98
13) 1,4-Dichlorobenzene	6.828	146	472812	37.916	ng	99
14) 1,2-Dichlorobenzene	6.975	146	436032	36.935	ng	97
15) Benzyl Alcohol	6.957	79	351040	40.337	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.081	45	524518	38.654	ng	97
17) 2-Methylphenol	7.063	107	345119	40.612	ng	97
18) Hexachloroethane	7.316	117	162292	37.103	ng	94
19) n-Nitroso-di-n-propyla...	7.228	70	294841	39.715	ng	96
20) 3+4-Methylphenols	7.216	107	433947	39.292	ng	92
22) Acetophenone	7.216	105	603475	37.477	ng	# 97
24) Nitrobenzene	7.393	77	430430	36.608	ng	96
25) Isophorone	7.628	82	803835	38.422	ng	97
26) 2-Nitrophenol	7.704	139	224890	44.890	ng	97
27) 2,4-Dimethylphenol	7.745	122	348673	52.878	ng	95
28) bis(2-Chloroethoxy)met...	7.840	93	496230	39.213	ng	97
29) 2,4-Dichlorophenol	7.945	162	366254	41.180	ng	98
30) 1,2,4-Trichlorobenzene	8.028	180	403318	37.039	ng	100
31) Naphthalene	8.110	128	1306143	37.901	ng	98
32) Benzoic acid	7.881	122	258186	42.963	ng	99
33) 4-Chloroaniline	8.157	127	281700	24.048	ng	98
34) Hexachlorobutadiene	8.222	225	229672	36.052	ng	99
35) Caprolactam	8.534	113	105041m	42.247	ng	
36) 4-Chloro-3-methylphenol	8.645	107	385995	40.015	ng	98
37) 2-Methylnaphthalene	8.798	142	821414	37.997	ng	98
38) 1-Methylnaphthalene	8.898	142	785945	37.269	ng	94
40) 1,2,4,5-Tetrachloroben...	8.969	216	392956	39.305	ng	98
41) Hexachlorocyclopentadiene	8.951	237	386205	100.393	ng	96
43) 2,4,6-Trichlorophenol	9.081	196	264409	40.746	ng	94

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120123\
 Data File : BF136332.D
 Acq On : 01 Dec 2023 12:08
 Operator : CG\JU
 Sample : PB157138BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB157138BS

Quant Time: Dec 01 13:16:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 03:04:20 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/04/2023
 Supervised By :mohammad ahmed 12/05/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.128	196	276311	39.018	ng	99	12/04/2023
46) 1,1'-Biphenyl	9.269	154	1060605	38.285	ng	99	Supervised By :mohammad
47) 2-Chloronaphthalene	9.292	162	803918	37.368	ng	98	ahmed
48) 2-Nitroaniline	9.392	65	225322	40.115	ng	# 84	
49) Acenaphthylene	9.710	152	1246747	40.609	ng	100	
50) Dimethylphthalate	9.575	163	926871	39.037	ng	99	
51) 2,6-Dinitrotoluene	9.634	165	196969	40.607	ng	99	12/05/2023
52) Acenaphthene	9.881	154	736039	37.937	ng	98	
53) 3-Nitroaniline	9.804	138	152525	31.690	ng	92	
54) 2,4-Dinitrophenol	9.916	184	192406	85.288	ng	# 87	
55) Dibenzofuran	10.057	168	1134050	39.281	ng	96	
56) 4-Nitrophenol	9.969	139	266935	73.024	ng	93	
57) 2,4-Dinitrotoluene	10.039	165	249344	38.483	ng	91	
58) Fluorene	10.398	166	849552	37.515	ng	99	
59) 2,3,4,6-Tetrachlorophenol	10.169	232	226600	40.217	ng	99	
60) Diethylphthalate	10.275	149	888217	37.397	ng	99	
61) 4-Chlorophenyl-phenyle...	10.392	204	415220	37.756	ng	99	
62) 4-Nitroaniline	10.422	138	160308	34.027	ng	95	
63) Azobenzene	10.551	77	818428	37.169	ng	96	
65) 4,6-Dinitro-2-methylph...	10.451	198	127219	43.281	ng	84	
66) n-Nitrosodiphenylamine	10.510	169	742811	43.123	ng	99	
67) 4-Bromophenyl-phenylether	10.881	248	261383	41.682	ng	98	
68) Hexachlorobenzene	10.945	284	294664	40.600	ng	95	
69) Atrazine	11.045	200	204750	40.070	ng	98	
70) Pentachlorophenol	11.139	266	263387	64.202	ng	98	
71) Phenanthrene	11.363	178	1236491	40.929	ng	99	
72) Anthracene	11.416	178	1229127	40.571	ng	99	
73) Carbazole	11.569	167	947114	36.757	ng	99	
74) Di-n-butylphthalate	11.904	149	1223653	37.955	ng	98	
75) Fluoranthene	12.557	202	1038900	33.875	ng	97	
77) Benzidine	12.680	184	94054	23.860	ng	99	
78) Pyrene	12.786	202	1013296	38.584	ng	97	
80) Butylbenzylphthalate	13.410	149	332815	40.719	ng	97	
81) Benzo(a)anthracene	13.980	228	702269	39.428	ng	98	
82) 3,3'-Dichlorobenzidine	13.945	252	205299	47.883	ng	97	
83) Chrysene	14.016	228	673921	39.187	ng	98	
84) Bis(2-ethylhexyl)phtha...	13.974	149	430367	43.874	ng	99	
85) Di-n-octyl phthalate	14.592	149	780485	48.884	ng	100	
87) Indeno(1,2,3-cd)pyrene	16.992	276	851352	44.023	ng	99	
88) Benzo(b)fluoranthene	15.039	252	733295	40.184	ng	99	
89) Benzo(k)fluoranthene	15.068	252	721501	41.629	ng	98	
90) Benzo(a)pyrene	15.410	252	647045	43.374	ng	97	
91) Dibenzo(a,h)anthracene	17.010	278	705136	45.347	ng	99	
92) Benzo(g,h,i)perylene	17.451	276	679502	43.578	ng	99	

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

Manual Integrations APPROVED

Quant Time: Dec 01 13:16:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 28 03:04:20 2023
Response via : Initial Calibration

