

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121223\
 Data File : BF136510.D
 Acq On : 12 Dec 2023 11:55
 Operator : CG\JU
 Sample : PB157670BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB157670BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/14/2023
 Supervised By :mohammad ahmed 12/14/2023

Quant Time: Dec 12 12:22:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 06 09:40:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	130641	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	550672	20.000	ng	0.00
39) Acenaphthene-d10	9.833	164	296630	20.000	ng	0.00
64) Phenanthrene-d10	11.328	188	562577	20.000	ng	0.00
76) Chrysene-d12	13.980	240	281029	20.000	ng	0.00
86) Perylene-d12	15.457	264	259818	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	996620	112.663	ng	0.02
7) Phenol-d6	6.451	99	1253496	113.574	ng	0.00
23) Nitrobenzene-d5	7.363	82	764087	75.012	ng	0.00
42) 2,4,6-Tribromophenol	10.633	330	434138	118.699	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1628748	74.923	ng	0.00
79) Terphenyl-d14	12.922	244	1780158	83.491	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.657	88	117442	33.566	ng	98
3) Pyridine	3.416	79	269946	31.858	ng	98
4) n-Nitrosodimethylamine	3.387	42	142014	38.577	ng	100
6) Aniline	6.463	93	397049	35.597	ng	# 10
8) 2-Chlorophenol	6.587	128	400871	41.510	ng	98
9) Benzaldehyde	6.351	77	227144	40.943	ng	95
10) Phenol	6.463	94	462939	39.419	ng	82
11) bis(2-Chloroethyl)ether	6.540	93	349961	39.978	ng	98
12) 1,3-Dichlorobenzene	6.740	146	402338	37.046	ng	98
13) 1,4-Dichlorobenzene	6.816	146	414581	37.631	ng	99
14) 1,2-Dichlorobenzene	6.969	146	395635	38.189	ng	97
15) Benzyl Alcohol	6.945	79	316274	42.794	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.075	45	425737	37.814	ng	94
17) 2-Methylphenol	7.063	107	322247	41.303	ng	94
18) Hexachloroethane	7.304	117	144919	37.676	ng	98
19) n-Nitroso-di-n-propyla...	7.216	70	254984	40.050	ng	99
20) 3+4-Methylphenols	7.216	107	393839	40.668	ng	# 85
22) Acetophenone	7.210	105	542795	39.917	ng	97
24) Nitrobenzene	7.381	77	382645	37.342	ng	98
25) Isophorone	7.622	82	722902	39.502	ng	97
26) 2-Nitrophenol	7.698	139	211511	41.986	ng	92
27) 2,4-Dimethylphenol	7.739	122	358696	57.745	ng	99
28) bis(2-Chloroethoxy)met...	7.834	93	458046	40.732	ng	99
29) 2,4-Dichlorophenol	7.939	162	339215	41.372	ng	99
30) 1,2,4-Trichlorobenzene	8.016	180	365927	38.274	ng	97
31) Naphthalene	8.098	128	1162157	38.270	ng	99
32) Benzoic acid	7.881	122	270000	43.816	ng	96
33) 4-Chloroaniline	8.151	127	304212	28.100	ng	99
34) Hexachlorobutadiene	8.210	225	209000	36.886	ng	98
35) Caprolactam	8.534	113	114346m	45.286	ng	
36) 4-Chloro-3-methylphenol	8.639	107	364794	41.903	ng	97
37) 2-Methylnaphthalene	8.792	142	755766	39.120	ng	96
38) 1-Methylnaphthalene	8.892	142	717939	37.871	ng	98
40) 1,2,4,5-Tetrachloroben...	8.957	216	365003	39.551	ng	99
41) Hexachlorocyclopentadiene	8.945	237	375487	109.825	ng	99
43) 2,4,6-Trichlorophenol	9.075	196	251442	41.844	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.122	196	272508	40.633	ng	96
46) 1,1'-Biphenyl	9.257	154	1004834	39.811	ng	98
47) 2-Chloronaphthalene	9.281	162	749688	38.162	ng	97
48) 2-Nitroaniline	9.381	65	216237	41.041	ng	95
49) Acenaphthylene	9.698	152	1130994	40.415	ng	99
50) Dimethylphthalate	9.569	163	906948	41.529	ng	99
51) 2,6-Dinitrotoluene	9.628	165	197302	40.158	ng	98
52) Acenaphthene	9.869	154	710512	39.621	ng	97
53) 3-Nitroaniline	9.798	138	168187	33.522	ng	99
54) 2,4-Dinitrophenol	9.910	184	219837	88.498	ng	98
55) Dibenzofuran	10.045	168	1043374	39.827	ng	99
56) 4-Nitrophenol	9.975	139	303591	80.719	ng	95
57) 2,4-Dinitrotoluene	10.033	165	267473	41.153	ng	95
58) Fluorene	10.386	166	827912	39.369	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.163	232	235229	43.787	ng	99
60) Diethylphthalate	10.269	149	873319	40.667	ng	99
61) 4-Chlorophenyl-phenyle...	10.381	204	412125	40.200	ng	95
62) 4-Nitroaniline	10.422	138	205263	41.203	ng	98
63) Azobenzene	10.539	77	768450	39.314	ng	96
65) 4,6-Dinitro-2-methylph...	10.445	198	152294	43.711	ng	94
66) n-Nitrosodiphenylamine	10.504	169	748784	41.479	ng	98
67) 4-Bromophenyl-phenylether	10.869	248	267508	40.898	ng	# 93
68) Hexachlorobenzene	10.933	284	302730	40.045	ng	99
69) Atrazine	11.039	200	310428	60.250	ng	98
70) Pentachlorophenol	11.133	266	307513	72.075	ng	98
71) Phenanthrene	11.357	178	1266727	40.966	ng	99
72) Anthracene	11.404	178	1273354	40.885	ng	99
73) Carbazole	11.563	167	1079480	39.160	ng	100
74) Di-n-butylphthalate	11.892	149	1364606	40.382	ng	99
75) Fluoranthene	12.545	202	1207194	37.254	ng	99
77) Benzidine	12.669	184	316415	45.309	ng	99
78) Pyrene	12.774	202	1187855	43.159	ng	99
80) Butylbenzylphthalate	13.398	149	422946	43.168	ng	92
81) Benzo(a)anthracene	13.969	228	810290	41.393	ng	99
82) 3,3'-Dichlorobenzidine	13.933	252	250805	45.696	ng	100
83) Chrysene	14.010	228	771973	40.083	ng	99
84) Bis(2-ethylhexyl)phtha...	13.963	149	507090	42.219	ng	99
85) Di-n-octyl phthalate	14.580	149	756769	42.655	ng	99
87) Indeno(1,2,3-cd)pyrene	16.962	276	851296	47.163	ng	99
88) Benzo(b)fluoranthene	15.027	252	725994	40.868	ng	100
89) Benzo(k)fluoranthene	15.057	252	695082	41.519	ng	99
90) Benzo(a)pyrene	15.398	252	631378	43.037	ng	99
91) Dibenzo(a,h)anthracene	16.992	278	710269	48.032	ng	100
92) Benzo(g,h,i)perylene	17.427	276	713709	47.023	ng	100

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

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