

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110824\
 Data File : BF140288.D
 Acq On : 08 Nov 2024 10:28
 Operator : RC/JU
 Sample : PB164750BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164750BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/10/2024
 Supervised By :mohammad ahmed 11/11/2024

Quant Time: Nov 08 11:05:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 05 16:47:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	195639	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	765547	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	473510	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	882492	20.000	ng	0.00
76) Chrysene-d12	14.063	240	554441	20.000	ng	0.00
86) Perylene-d12	15.562	264	455207	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1350884	114.430	ng	0.02
7) Phenol-d6	6.510	99	1707980	112.481	ng	0.00
23) Nitrobenzene-d5	7.445	82	1176601	76.414	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	676037	144.440	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	2226512	75.084	ng	0.00
79) Terphenyl-d14	12.998	244	2859948	84.499	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.751	88	162998	30.032	ng	96
3) Pyridine	3.498	79	445970	33.675	ng	95
4) n-Nitrosodimethylamine	3.457	42	274935	35.958	ng	90
6) Aniline	6.539	93	648715	47.287	ng	92
8) 2-Chlorophenol	6.663	128	528701	42.803	ng	95
9) Benzaldehyde	6.428	77	67892	7.258	ng	94
10) Phenol	6.528	94	668118	41.429	ng	94
11) bis(2-Chloroethyl)ether	6.616	93	503595	41.107	ng	96
12) 1,3-Dichlorobenzene	6.822	146	559655	38.664	ng	98
13) 1,4-Dichlorobenzene	6.898	146	567610	38.881	ng	98
14) 1,2-Dichlorobenzene	7.045	146	546411	40.074	ng	99
15) Benzyl Alcohol	7.016	79	499648	42.685	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.151	45	782413	39.455	ng	98
17) 2-Methylphenol	7.128	107	434126	42.132	ng	99
18) Hexachloroethane	7.392	117	212125	39.008	ng	98
19) n-Nitroso-di-n-propyla...	7.292	70	393754	40.647	ng	96
20) 3+4-Methylphenols	7.281	107	533406	41.218	ng	97
22) Acetophenone	7.286	105	719910	37.929	ng	97
24) Nitrobenzene	7.463	77	589350	36.477	ng	96
25) Isophorone	7.698	82	1076954	40.620	ng	99
26) 2-Nitrophenol	7.775	139	281854	43.855	ng	96
27) 2,4-Dimethylphenol	7.810	122	436796	49.957	ng	96
28) bis(2-Chloroethoxy)met...	7.904	93	630073	39.712	ng	99
29) 2,4-Dichlorophenol	8.016	162	459059	42.440	ng	99
30) 1,2,4-Trichlorobenzene	8.098	180	490581	37.849	ng	99
31) Naphthalene	8.181	128	1536642	38.871	ng	100
32) Benzoic acid	7.939	122	318810	44.431	ng	96
33) 4-Chloroaniline	8.228	127	515740	39.629	ng	97
34) Hexachlorobutadiene	8.298	225	337982	38.524	ng	99
35) Caprolactam	8.610	113	131452m	39.949	ng	
36) 4-Chloro-3-methylphenol	8.710	107	518752	43.075	ng	97
37) 2-Methylnaphthalene	8.869	142	1051398	40.752	ng	99
38) 1-Methylnaphthalene	8.969	142	974624	38.541	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	528119	38.098	ng	99
41) Hexachlorocyclopentadiene	9.028	237	641602	165.176	ng	98
43) 2,4,6-Trichlorophenol	9.151	196	369763	43.045	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	371986	41.185	ng	98
46) 1,1'-Biphenyl	9.339	154	1294320	37.817	ng	100
47) 2-Chloronaphthalene	9.363	162	970176	37.662	ng	100
48) 2-Nitroaniline	9.463	65	345821	40.142	ng	92
49) Acenaphthylene	9.780	152	1541359	42.263	ng	99
50) Dimethylphthalate	9.645	163	1234962	42.349	ng	100
51) 2,6-Dinitrotoluene	9.704	165	282787	41.254	ng	96
52) Acenaphthene	9.957	154	1173915	45.381	ng	99
53) 3-Nitroaniline	9.875	138	239845	37.159	ng	95
54) 2,4-Dinitrophenol	9.986	184	324620	95.234	ng	87
55) Dibenzofuran	10.127	168	1452145	40.145	ng	98
56) 4-Nitrophenol	10.039	139	426153	93.901	ng	98
57) 2,4-Dinitrotoluene	10.116	165	380332	42.757	ng	97
58) Fluorene	10.475	166	1149420	39.919	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.245	232	350955	47.432	ng	95
60) Diethylphthalate	10.351	149	1226888	42.337	ng	99
61) 4-Chlorophenyl-phenyle...	10.463	204	585618	40.355	ng	96
62) 4-Nitroaniline	10.498	138	286478	44.663	ng	95
63) Azobenzene	10.622	77	1209988	40.279	ng	98
65) 4,6-Dinitro-2-methylph...	10.527	198	233285	50.332	ng	91
66) n-Nitrosodiphenylamine	10.586	169	1025499	40.403	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	399731	42.064	ng	96
68) Hexachlorobenzene	11.022	284	447314	41.295	ng	95
69) Atrazine	11.110	200	332139	47.829	ng	98
70) Pentachlorophenol	11.216	266	541719	93.134	ng	98
71) Phenanthrene	11.439	178	1745451	41.444	ng	100
72) Anthracene	11.492	178	1767910	42.809	ng	100
73) Carbazole	11.645	167	1547770	41.017	ng	100
74) Di-n-butylphthalate	11.969	149	1902404	42.591	ng	99
75) Fluoranthene	12.627	202	1842695	41.844	ng	100
77) Benzidine	12.745	184	485628	44.798	ng	99
78) Pyrene	12.857	202	1874116	40.939	ng	100
80) Butylbenzylphthalate	13.474	149	750677	45.539	ng	94
81) Benzo(a)anthracene	14.051	228	1497295	42.081	ng	100
82) 3,3'-Dichlorobenzidine	14.015	252	449354	40.215	ng	99
83) Chrysene	14.092	228	1402669	42.096	ng	100
84) Bis(2-ethylhexyl)phtha...	14.033	149	1033489	44.733	ng	99
85) Di-n-octyl phthalate	14.662	149	1560464	44.874	ng	96
87) Indeno(1,2,3-cd)pyrene	17.080	276	1169241	43.758	ng	99
88) Benzo(b)fluoranthene	15.121	252	1214774	44.118	ng	100
89) Benzo(k)fluoranthene	15.157	252	1073254	42.478	ng	99
90) Benzo(a)pyrene	15.498	252	1045593	46.058	ng	100
91) Dibenzo(a,h)anthracene	17.098	278	946410	43.051	ng	99
92) Benzo(g,h,i)perylene	17.539	276	875549	39.042	ng	99

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

