

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070823\
 Data File : BF134163.D
 Acq On : 09 Jul 2023 00:01
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 07/09/2023

Supervised By :mohammad ahmed 07/09/2023

Quant Time: Jul 09 00:42:22 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jun 30 18:35:29 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	167584	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	623095	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	316872	20.000	ng	0.00
64) Phenanthrene-d10	11.380	188	608217	20.000	ng	0.00
76) Chrysene-d12	14.045	240	349490	20.000	ng	0.00
86) Perylene-d12	15.545	264	297883	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	781101	77.967	ng	0.00
7) Phenol-d6	6.492	99	975555	79.181	ng	0.00
23) Nitrobenzene-d5	7.416	82	926408	80.146	ng	0.00
42) 2,4,6-Tribromophenol	10.680	330	315414	79.391	ng	0.00
45) 2-Fluorobiphenyl	9.210	172	1647405	75.587	ng	0.00
79) Terphenyl-d14	12.980	244	1724998	79.437	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.640	88	147540	35.716	ng	97
3) Pyridine	3.416	79	420447	39.075	ng	98
4) n-Nitrosodimethylamine	3.393	42	228576	37.195	ng	99
6) Aniline	6.522	93	619015	41.288	ng	99
8) 2-Chlorophenol	6.639	128	380863	37.450	ng	94
9) Benzaldehyde	6.404	77	279224	41.714	ng	99
10) Phenol	6.504	94	507663	39.189	ng	95
11) bis(2-Chloroethyl)ether	6.592	93	388741	40.761	ng	98
12) 1,3-Dichlorobenzene	6.786	146	412649	38.059	ng	98
13) 1,4-Dichlorobenzene	6.863	146	419560	38.312	ng	98
14) 1,2-Dichlorobenzene	7.022	146	390881	38.112	ng	98
15) Benzyl Alcohol	6.992	79	367324	38.674	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.128	45	653297	38.720	ng	98
17) 2-Methylphenol	7.104	107	363461	39.910	ng	99
18) Hexachloroethane	7.357	117	164574	40.530	ng	97
19) n-Nitroso-di-n-propyla...	7.269	70	306745	39.259	ng	97
20) 3+4-Methylphenols	7.257	107	427368	38.682	ng	# 86
22) Acetophenone	7.263	105	552178	38.966	ng	97
24) Nitrobenzene	7.433	77	476416	40.787	ng	96
25) Isophorone	7.669	82	822377	39.146	ng	98
26) 2-Nitrophenol	7.745	139	240505	42.921	ng	98
27) 2,4-Dimethylphenol	7.786	122	355361	40.064	ng	98
28) bis(2-Chloroethoxy)met...	7.880	93	494129	40.622	ng	98
29) 2,4-Dichlorophenol	7.986	162	344413	39.158	ng	96
30) 1,2,4-Trichlorobenzene	8.069	180	364080	40.117	ng	98
31) Naphthalene	8.151	128	1146459	38.092	ng	99
32) Benzoic acid	7.922	122	266680m	40.124	ng	
33) 4-Chloroaniline	8.204	127	503755	39.128	ng	98
34) Hexachlorobutadiene	8.263	225	233394	40.768	ng	100
35) Caprolactam	8.592	113	113809	39.298	ng	96
36) 4-Chloro-3-methylphenol	8.686	107	393289	39.803	ng	100
37) 2-Methylnaphthalene	8.839	142	822305	38.636	ng	99
38) 1-Methylnaphthalene	8.939	142	762573	38.541	ng	99
40) 1,2,4,5-Tetrachloroben...	9.010	216	377970	40.262	ng	100
41) Hexachlorocyclopentadiene	8.992	237	145574	32.686	ng	99
43) 2,4,6-Trichlorophenol	9.122	196	249356	39.019	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.163	196	292055	38.851	ng	97
46) 1,1'-Biphenyl	9.310	154	937800	36.895	ng	97
47) 2-Chloronaphthalene	9.333	162	700990	37.692	ng	97
48) 2-Nitroaniline	9.433	65	262155	38.678	ng	99
49) Acenaphthylene	9.751	152	1189081	37.428	ng	99
50) Dimethylphthalate	9.616	163	903039	39.259	ng	100
51) 2,6-Dinitrotoluene	9.680	165	199514	40.272	ng	96
52) Acenaphthene	9.922	154	709629	38.495	ng	99
53) 3-Nitroaniline	9.851	138	228928	38.518	ng	99
54) 2,4-Dinitrophenol	9.957	184	94923	35.439	ng	98
55) Dibenzofuran	10.092	168	1076761	37.374	ng	98
56) 4-Nitrophenol	10.010	139	149464	35.952	ng	95
57) 2,4-Dinitrotoluene	10.086	165	252283	38.560	ng	91
58) Fluorene	10.439	166	824407	36.781	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.216	232	229147	38.645	ng	97
60) Diethylphthalate	10.310	149	942241	39.318	ng	98
61) 4-Chlorophenyl-phenyle...	10.427	204	411409	38.088	ng	96
62) 4-Nitroaniline	10.469	138	227255	37.941	ng	99
63) Azobenzene	10.592	77	878111	38.737	ng	97
65) 4,6-Dinitro-2-methylph...	10.498	198	141051	42.537	ng	96
66) n-Nitrosodiphenylamine	10.551	169	751856	38.690	ng	97
67) 4-Bromophenyl-phenylether	10.921	248	262753	40.645	ng	95
68) Hexachlorobenzene	10.986	284	278513	40.577	ng	98
69) Atrazine	11.080	200	222579	41.408	ng	98
70) Pentachlorophenol	11.186	266	155283	37.837	ng	98
71) Phenanthrene	11.410	178	1146073	37.431	ng	100
72) Anthracene	11.463	178	1212618	38.076	ng	100
73) Carbazole	11.616	167	1101714	37.795	ng	100
74) Di-n-butylphthalate	11.945	149	1443683	41.647	ng	99
75) Fluoranthene	12.604	202	1249129	37.736	ng	98
77) Benzidine	12.727	184	365740	43.981	ng	99
78) Pyrene	12.833	202	1163167	35.762	ng	100
80) Butylbenzylphthalate	13.457	149	560325	45.935	ng	95
81) Benzo(a)anthracene	14.033	228	929078	38.332	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	304657	39.674	ng	# 98
83) Chrysene	14.068	228	903698	38.495	ng	100
84) Bis(2-ethylhexyl)phtha...	14.015	149	736053	52.513	ng	100
85) Di-n-octyl phthalate	14.639	149	1082379	52.554	ng	99
87) Indeno(1,2,3-cd)pyrene	17.098	276	848361	41.043	ng	99
88) Benzo(b)fluoranthene	15.104	252	753310	43.008	ng	100
89) Benzo(k)fluoranthene	15.133	252	756413	42.415	ng	99
90) Benzo(a)pyrene	15.480	252	670701	39.598	ng	99
91) Dibenzo(a,h)anthracene	17.115	278	704644	41.737	ng	98
92) Benzo(g,h,i)perylene	17.568	276	645317	37.065	ng	98

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

