

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080122\
 Data File : BM036035.D
 Acq On : 01 Aug 2022 17:29
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM080122

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/02/2022
 Supervised By : Jagrut Upadhyay 08/02/2022

Quant Time: Aug 01 18:53:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	257484	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1097771	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	660682	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	1344660	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1358307	20.000	ng	0.00
86) Perylene-d12	23.791	264	1381754	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1173164	73.869	ng	0.00
7) Phenol-d6	7.039	99	1486361	73.552	ng	0.00
23) Nitrobenzene-d5	9.039	82	1395023	71.134	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	560300	72.291	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	3096100	69.191	ng	0.00
79) Terphenyl-d14	19.874	244	4751134	68.993	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.299	88	234172	36.591	ng	# 92
3) Pyridine	3.710	79	629687	36.693	ng	89
4) n-Nitrosodimethylamine	3.622	42	256291	36.343	ng	# 62
6) Aniline	7.198	93	826186	36.919	ng	96
8) 2-Chlorophenol	7.434	128	624585	36.885	ng	95
9) Benzaldehyde	7.010	77	356095	33.285	ng	94
10) Phenol	7.063	94	751093	36.913	ng	91
11) bis(2-Chloroethyl)ether	7.298	93	608046	36.276	ng	95
12) 1,3-Dichlorobenzene	7.757	146	687867	36.658	ng	97
13) 1,4-Dichlorobenzene	7.904	146	699990	36.580	ng	99
14) 1,2-Dichlorobenzene	8.222	146	680668	36.865	ng	99
15) Benzyl Alcohol	8.116	79	497952	36.720	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.410	45	805603	36.144	ng	96
17) 2-Methylphenol	8.316	107	497070	36.904	ng	95
18) Hexachloroethane	8.951	117	254135	36.892	ng	97
19) n-Nitroso-di-n-propyla...	8.686	70	455256	37.088	ng	# 87
20) 3+4-Methylphenols	8.651	107	675605	36.919	ng	95
22) Acetophenone	8.698	105	915929	35.247	ng	# 93
24) Nitrobenzene	9.081	77	656953	35.634	ng	97
25) Isophorone	9.610	82	1122552	35.186	ng	98
26) 2-Nitrophenol	9.792	139	315703	36.306	ng	91
27) 2,4-Dimethylphenol	9.851	122	478090	35.492	ng	97
28) bis(2-Chloroethoxy)met...	10.092	93	789680	34.743	ng	100
29) 2,4-Dichlorophenol	10.322	162	544129	35.920	ng	99
30) 1,2,4-Trichlorobenzene	10.539	180	604101	35.190	ng	100
31) Naphthalene	10.727	128	1910371	35.002	ng	99
32) Benzoic acid	9.986	122	394438	36.940	ng	93
33) 4-Chloroaniline	10.839	127	781794	35.543	ng	96
34) Hexachlorobutadiene	11.010	225	357420	35.432	ng	96
35) Caprolactam	11.633	113	173405	37.164	ng	90
36) 4-Chloro-3-methylphenol	11.963	107	568865	35.727	ng	98
37) 2-Methylnaphthalene	12.339	142	1279063	35.217	ng	97
38) 1-Methylnaphthalene	12.557	142	1250440	35.157	ng	98
40) 1,2,4,5-Tetrachloroben...	12.704	216	618974	34.557	ng	99
41) Hexachlorocyclopentadiene	12.680	237	335797	35.397	ng	100
43) 2,4,6-Trichlorophenol	12.945	196	432315	35.470	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.016	196	459507	35.715	ng	98
46) 1,1'-Biphenyl	13.345	154	1658776	34.451	ng	99
47) 2-Chloronaphthalene	13.386	162	1270548	34.478	ng	99
48) 2-Nitroaniline	13.598	65	363766	36.222	ng	92
49) Acenaphthylene	14.233	152	2004709	35.193	ng	99
50) Dimethylphthalate	13.974	163	1565288	34.839	ng	99
51) 2,6-Dinitrotoluene	14.092	165	327430	36.322	ng	93
52) Acenaphthene	14.574	154	1302173m	34.946	ng	
53) 3-Nitroaniline	14.421	138	385835	36.685	ng	93
54) 2,4-Dinitrophenol	14.627	184	176260	36.949	ng	87
55) Dibenzofuran	14.904	168	1943949	34.837	ng	98
56) 4-Nitrophenol	14.721	139	315532	37.490	ng	89
57) 2,4-Dinitrotoluene	14.874	165	436384	36.933	ng	99
58) Fluorene	15.551	166	1532110	34.807	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.133	232	389284	35.718	ng	97
60) Diethylphthalate	15.333	149	1623164	35.155	ng	99
61) 4-Chlorophenyl-phenyle...	15.551	204	764954	34.852	ng	96
62) 4-Nitroaniline	15.580	138	408455	37.751	ng	89
63) Azobenzene	15.845	77	1553867	35.401	ng	93
65) 4,6-Dinitro-2-methylph...	15.633	198	252197	36.681	ng	94
66) n-Nitrosodiphenylamine	15.762	169	1309812	34.506	ng	100
67) 4-Bromophenyl-phenylether	16.445	248	463270	34.444	ng	95
68) Hexachlorobenzene	16.551	284	533855	34.593	ng	92
69) Atrazine	16.715	200	405038	38.134	ng	98
70) Pentachlorophenol	16.898	266	328133	36.305	ng	98
71) Phenanthrene	17.292	178	2378380	34.718	ng	99
72) Anthracene	17.380	178	2337747	35.248	ng	99
73) Carbazole	17.651	167	2406885	35.817	ng	99
74) Di-n-butylphthalate	18.221	149	2869696	35.773	ng	99
75) Fluoranthene	19.303	202	2752273	35.891	ng	98
77) Benzidine	19.492	184	802718	39.406	ng	99
78) Pyrene	19.668	202	2872561	34.104	ng	99
80) Butylbenzylphthalate	20.568	149	1280452	35.222	ng	99
81) Benzo(a)anthracene	21.415	228	2918758	34.887	ng	99
82) 3,3'-Dichlorobenzidine	21.350	252	729267	35.898	ng	99
83) Chrysene	21.468	228	2799507	35.201	ng	99
84) Bis(2-ethylhexyl)phtha...	21.350	149	1923127	35.065	ng	# 98
85) Di-n-octyl phthalate	22.262	149	3209836	35.861	ng	95
87) Indeno(1,2,3-cd)pyrene	26.256	276	3388742	34.899	ng	98
88) Benzo(b)fluoranthene	23.074	252	2829353	34.949	ng	99
89) Benzo(k)fluoranthene	23.121	252	2847190	34.821	ng	99
90) Benzo(a)pyrene	23.691	252	2379969	35.039	ng	99
91) Dibenzo(a,h)anthracene	26.273	278	2852757	35.098	ng	99
92) Benzo(g,h,i)perylene	27.009	276	2755615	35.020	ng	98

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

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