

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011422\
 Data File : BP008809.D
 Acq On : 14 Jan 2022 15:33
 Operator : CG/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/17/2022
 Supervised By : Jagrut Upadhyay 01/17/2022

Quant Time: Jan 14 16:09:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 14 16:06:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	324714	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1268154	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	801278	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	1513989	20.000	ng	0.00
76) Chrysene-d12	21.345	240	1282268	20.000	ng	0.00
86) Perylene-d12	23.762	264	1324554	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1481844	71.463	ng	0.00
7) Phenol-d6	7.034	99	1850923	66.493	ng	0.00
23) Nitrobenzene-d5	9.016	82	1616632	72.195	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	836729	61.896	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	4285965	75.200	ng	0.00
79) Terphenyl-d14	19.815	244	5647057	89.746	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.346	88	287676	36.785	ng	97
3) Pyridine	3.746	79	640785	34.832	ng	97
4) n-Nitrosodimethylamine	3.652	42	300435	36.651	ng	86
6) Aniline	7.181	93	1152817	32.967	ng	99
8) 2-Chlorophenol	7.422	128	804958	34.349	ng	98
9) Benzaldehyde	6.993	77	608779	32.314	ng	97
10) Phenol	7.057	94	938495	32.947	ng	97
11) bis(2-Chloroethyl)ether	7.281	93	740460	33.537	ng	99
12) 1,3-Dichlorobenzene	7.746	146	946089	35.735	ng	99
13) 1,4-Dichlorobenzene	7.893	146	953781	35.253	ng	99
14) 1,2-Dichlorobenzene	8.210	146	909882	34.681	ng	99
15) Benzyl Alcohol	8.098	79	657539	31.797	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.387	45	838064	33.302	ng	99
17) 2-Methylphenol	8.304	107	634527	32.312	ng	97
18) Hexachloroethane	8.940	117	329541	36.212	ng	97
19) n-Nitroso-di-n-propyla...	8.669	70	554602	30.686	ng	97
20) 3+4-Methylphenols	8.640	107	853439	31.322	ng	97
22) Acetophenone	8.681	105	1166327	34.856	ng	# 97
24) Nitrobenzene	9.057	77	815285	35.966	ng	96
25) Isophorone	9.592	82	1480764	33.623	ng	99
26) 2-Nitrophenol	9.769	139	437730	37.028	ng	95
27) 2,4-Dimethylphenol	9.840	122	739839	34.939	ng	99
28) bis(2-Chloroethoxy)met...	10.069	93	953186	34.712	ng	99
29) 2,4-Dichlorophenol	10.316	162	801116	35.292	ng	99
30) 1,2,4-Trichlorobenzene	10.522	180	914757	36.990	ng	99
31) Naphthalene	10.710	128	2437818	35.554	ng	100
32) Benzoic acid	10.010	122	451714	31.907	ng	99
33) 4-Chloroaniline	10.816	127	1024537	33.293	ng	98
34) Hexachlorobutadiene	10.998	225	574520	38.426	ng	99
35) Caprolactam	11.616	113	223897	27.756	ng	97
36) 4-Chloro-3-methylphenol	11.957	107	731183	31.802	ng	98
37) 2-Methylnaphthalene	12.322	142	1809895	33.550	ng	98
38) 1-Methylnaphthalene	12.539	142	1677705	33.485	ng	99
40) 1,2,4,5-Tetrachloroben...	12.692	216	1036962	38.932	ng	99
41) Hexachlorocyclopentadiene	12.675	237	576168	36.386	ng	100
43) 2,4,6-Trichlorophenol	12.933	196	657509	37.076	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.010	196	705277	36.180	ng	99
46) 1,1'-Biphenyl	13.333	154	2306359	36.924	ng	99
47) 2-Chloronaphthalene	13.369	162	1781799	37.166	ng	99
48) 2-Nitroaniline	13.575	65	408193	34.882	ng	97
49) Acenaphthylene	14.216	152	2711502	35.601	ng	100
50) Dimethylphthalate	13.957	163	2107391	33.095	ng	100
51) 2,6-Dinitrotoluene	14.075	165	459674	33.381	ng	96
52) Acenaphthene	14.563	154	1606275	35.212	ng	100
53) 3-Nitroaniline	14.404	138	449856	32.008	ng	98
54) 2,4-Dinitrophenol	14.616	184	192039	27.345	ng	92
55) Dibenzofuran	14.892	168	2605293	34.631	ng	98
56) 4-Nitrophenol	14.722	139	331480	28.324	ng	92
57) 2,4-Dinitrotoluene	14.863	165	600385	31.942	ng	93
58) Fluorene	15.545	166	2074767	34.261	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.127	232	581866m	33.216	ng	
60) Diethylphthalate	15.322	149	2013303	32.607	ng	99
61) 4-Chlorophenyl-phenyle...	15.539	204	1135262	34.941	ng	98
62) 4-Nitroaniline	15.569	138	437656	29.992	ng	94
63) Azobenzene	15.833	77	1713301	35.399	ng	94
65) 4,6-Dinitro-2-methylph...	15.633	198	290012	34.525	ng	94
66) n-Nitrosodiphenylamine	15.757	169	1769789	38.573	ng	99
67) 4-Bromophenyl-phenylether	16.439	248	696492	38.882	ng	99
68) Hexachlorobenzene	16.563	284	781640	37.504	ng	99
69) Atrazine	16.716	200	658418	35.353	ng	99
70) Pentachlorophenol	16.910	266	425371	32.694	ng	99
71) Phenanthrene	17.292	178	3036933	36.103	ng	99
72) Anthracene	17.386	178	3096901	36.050	ng	99
73) Carbazole	17.657	167	2615217	33.217	ng	99
74) Di-n-butylphthalate	18.210	149	3238807	35.930	ng	99
75) Fluoranthene	19.263	202	3311738	32.878	ng	97
77) Benzidine	19.439	184	1821932	33.207	ng	98
78) Pyrene	19.615	202	3337135	39.802	ng	98
80) Butylbenzylphthalate	20.492	149	1331469	38.839	ng	98
81) Benzo(a)anthracene	21.327	228	3081933	36.636	ng	98
82) 3,3'-Dichlorobenzidine	21.257	252	1356274	35.489	ng	99
83) Chrysene	21.380	228	2995191	36.910	ng	97
84) Bis(2-ethylhexyl)phtha...	21.257	149	2076607	41.489	ng	98
85) Di-n-octyl phthalate	22.186	149	3189101	35.898	ng	99
87) Indeno(1,2,3-cd)pyrene	26.297	276	3886762	36.723	ng	98
88) Benzo(b)fluoranthene	23.027	252	3131092	36.179	ng	99
89) Benzo(k)fluoranthene	23.074	252	2957276	36.511	ng	98
90) Benzo(a)pyrene	23.656	252	3011651	36.004	ng	99
91) Dibenzo(a,h)anthracene	26.315	278	3318888	37.112	ng	98
92) Benzo(g,h,i)perylene	27.074	276	3257097	36.691	ng	98

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

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