

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060624\
 Data File : BP020537.D
 Acq On : 06 Jun 2024 11:30
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/07/2024
 Supervised By :mohammad ahmed 06/08/2024

Quant Time: Jun 06 13:04:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 03:41:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	334983	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1298318	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	759733	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1298216	20.000	ng	-0.01
76) Chrysene-d12	21.639	240	1035158	20.000	ng	-0.03
86) Perylene-d12	25.004	264	1347135	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1485063	79.375	ng	0.00
7) Phenol-d6	6.934	99	1939523	73.509	ng	0.00
23) Nitrobenzene-d5	8.910	82	1889337	79.659	ng	-0.01
42) 2,4,6-Tribromophenol	15.904	330	685132	86.895	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	4296847	84.312	ng	0.00
79) Terphenyl-d14	19.933	244	4444852	71.481	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.322	88	293984	38.888	ng	96
3) Pyridine	3.711	79	803877m	37.821	ng	
4) n-Nitrosodimethylamine	3.628	42	362722	35.134	ng	96
6) Aniline	7.104	93	936953	35.045	ng	98
8) 2-Chlorophenol	7.328	128	826354	39.401	ng	97
9) Benzaldehyde	6.916	77	558027	32.971	ng	96
10) Phenol	6.957	94	981445	36.314	ng	99
11) bis(2-Chloroethyl)ether	7.193	93	816120	36.616	ng	99
12) 1,3-Dichlorobenzene	7.652	146	980184	39.667	ng	99
13) 1,4-Dichlorobenzene	7.799	146	983820	39.172	ng	99
14) 1,2-Dichlorobenzene	8.110	146	946789	39.062	ng	99
15) Benzyl Alcohol	7.993	79	689744	35.391	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.281	45	1094547	32.893	ng	98
17) 2-Methylphenol	8.199	107	673853	36.741	ng	100
18) Hexachloroethane	8.828	117	364974	39.798	ng	94
19) n-Nitroso-di-n-propyla...	8.563	70	624000	34.643	ng	96
20) 3+4-Methylphenols	8.528	107	933512	37.422	ng	99
22) Acetophenone	8.575	105	1251374	38.830	ng	# 98
24) Nitrobenzene	8.951	77	935447	38.862	ng	97
25) Isophorone	9.475	82	1714734	37.210	ng	99
26) 2-Nitrophenol	9.651	139	427755	45.166	ng	98
27) 2,4-Dimethylphenol	9.710	122	536904	38.565	ng	99
28) bis(2-Chloroethoxy)met...	9.951	93	1041679	36.777	ng	100
29) 2,4-Dichlorophenol	10.181	162	772795	41.376	ng	99
30) 1,2,4-Trichlorobenzene	10.398	180	909539	41.020	ng	98
31) Naphthalene	10.587	128	2624613	39.544	ng	100
32) Benzoic acid	9.857	122	556064	42.892	ng	99
33) 4-Chloroaniline	10.698	127	904462	35.227	ng	99
34) Hexachlorobutadiene	10.863	225	606294	43.408	ng	99
35) Caprolactam	11.487	113	225563	32.955	ng	91
36) 4-Chloro-3-methylphenol	11.810	107	798103	36.496	ng	99
37) 2-Methylnaphthalene	12.187	142	1757159	36.947	ng	98
38) 1-Methylnaphthalene	12.428	142	1721621	36.517	ng	99
40) 1,2,4,5-Tetrachloroben...	12.557	216	971153	44.180	ng	99
41) Hexachlorocyclopentadiene	12.534	237	404525	52.363	ng	99
43) 2,4,6-Trichlorophenol	12.804	196	614979	46.283	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.869	196	664162	43.327	ng	98
46) 1,1'-Biphenyl	13.210	154	2222819	41.457	ng	99
47) 2-Chloronaphthalene	13.257	162	1776603	41.681	ng	99
48) 2-Nitroaniline	13.463	65	470904	40.772	ng	98
49) Acenaphthylene	14.104	152	2529377	39.992	ng	100
50) Dimethylphthalate	13.857	163	1993434	37.224	ng	99
51) 2,6-Dinitrotoluene	13.969	165	452060	39.493	ng	97
52) Acenaphthene	14.457	154	1587855	39.800	ng	99
53) 3-Nitroaniline	14.304	138	411145	35.808	ng	95
54) 2,4-Dinitrophenol	14.492	184	188522	44.528	ng	95
55) Dibenzofuran	14.792	168	2411906	38.164	ng	99
56) 4-Nitrophenol	14.604	139	313333	35.278	ng	95
57) 2,4-Dinitrotoluene	14.751	165	564187	37.505	ng	96
58) Fluorene	15.457	166	1909149	37.630	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.028	232	523631	41.741	ng	97
60) Diethylphthalate	15.228	149	1907390	35.038	ng	100
61) 4-Chlorophenyl-phenyle...	15.457	204	1021974	39.035	ng	98
62) 4-Nitroaniline	15.481	138	374969	31.703	ng	97
63) Azobenzene	15.757	77	1751232	34.451	ng	96
65) 4,6-Dinitro-2-methylph...	15.539	198	266497	46.573	ng	93
66) n-Nitrosodiphenylamine	15.669	169	1568649	43.949	ng	99
67) 4-Bromophenyl-phenylether	16.363	248	600930	45.246	ng	98
68) Hexachlorobenzene	16.475	284	688814	44.067	ng	99
69) Atrazine	16.645	200	463132	36.882	ng	99
70) Pentachlorophenol	16.833	266	421326	44.380	ng	98
71) Phenanthrene	17.239	178	2585464	40.514	ng	99
72) Anthracene	17.327	178	2501571	39.948	ng	100
73) Carbazole	17.610	167	2208703	36.464	ng	100
74) Di-n-butylphthalate	18.210	149	2648029	36.100	ng	99
75) Fluoranthene	19.322	202	2701942	35.788	ng	99
77) Benzidine	19.533	184	763803m	24.143	ng	
78) Pyrene	19.698	202	2735197	35.128	ng	100
80) Butylbenzylphthalate	20.663	149	1098684	35.695	ng	95
81) Benzo(a)anthracene	21.621	228	2643555	38.785	ng	99
82) 3,3'-Dichlorobenzidine	21.545	252	981420	48.438	ng	99
83) Chrysene	21.692	228	2565467	39.945	ng	99
84) Bis(2-ethylhexyl)phtha...	21.580	149	1696949	38.294	ng	98
85) Di-n-octyl phthalate	22.874	149	3142184	48.786	ng	97
87) Indeno(1,2,3-cd)pyrene	28.815	276	3805145	46.450	ng	99
88) Benzo(b)fluoranthene	23.945	252	3065244	36.524	ng	98
89) Benzo(k)fluoranthene	24.015	252	3026015	36.969	ng	99
90) Benzo(a)pyrene	24.845	252	2792824	40.182	ng	100
91) Dibenzo(a,h)anthracene	28.909	278	3150463	46.108	ng	97
92) Benzo(g,h,i)perylene	30.009	276	3140773	45.988	ng	98

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

