

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052423\
 Data File : BP015274.D
 Acq On : 24 May 2023 16:05
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 05/25/2023
 Supervised By :Jagrut Upadhyay 05/25/2023

Quant Time: May 25 01:22:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 25 01:19:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.111	152	167818	20.000	ng	0.00
21) Naphthalene-d8	10.940	136	655718	20.000	ng	0.00
39) Acenaphthene-d10	14.751	164	385834	20.000	ng	0.00
64) Phenanthrene-d10	17.504	188	797314	20.000	ng	0.00
76) Chrysene-d12	21.581	240	752905	20.000	ng	0.00
86) Perylene-d12	24.139	264	857805	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.634	112	111058	10.966	ng	0.00
7) Phenol-d6	7.258	99	142070	10.704	ng	0.00
23) Nitrobenzene-d5	9.287	82	133811	9.964	ng	0.00
42) 2,4,6-Tribromophenol	16.240	330	48429	9.236	ng	0.00
45) 2-Fluorobiphenyl	13.375	172	322714	11.229	ng	0.00
79) Terphenyl-d14	20.039	244	442107	11.045	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.452	88	25652	5.719	ng	97
3) Pyridine	3.887	79	55292	4.823	ng	# 88
4) n-Nitrosodimethylamine	3.787	42	28387	5.306	ng	# 97
6) Aniline	7.428	93	87311	5.205	ng	99
8) 2-Chlorophenol	7.669	128	57241	5.409	ng	97
10) Phenol	7.287	94	71503	5.379	ng	97
11) bis(2-Chloroethyl)ether	7.522	93	62103	5.574	ng	99
12) 1,3-Dichlorobenzene	7.999	146	67591	5.740	ng	99
13) 1,4-Dichlorobenzene	8.146	146	66644	5.622	ng	99
14) 1,2-Dichlorobenzene	8.469	146	63931	5.653	ng	98
15) Benzyl Alcohol	8.352	79	43512	4.697	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.640	45	80564	5.607	ng	98
17) 2-Methylphenol	8.558	107	48322	5.108	ng	97
18) Hexachloroethane	9.199	117	25265	5.655	ng	92
19) n-Nitroso-di-n-propyla...	8.922	70	44346	5.221	ng	100
20) 3+4-Methylphenols	8.887	107	63827	4.995	ng	97
22) Acetophenone	8.946	105	88786	5.213	ng	# 96
24) Nitrobenzene	9.328	77	67220	5.027	ng	99
25) Isophorone	9.858	82	117958	4.894	ng	98
26) 2-Nitrophenol	10.052	139	25400	4.328	ng	99
27) 2,4-Dimethylphenol	10.099	122	49366	5.017	ng	95
28) bis(2-Chloroethoxy)met...	10.340	93	77512	5.218	ng	98
29) 2,4-Dichlorophenol	10.587	162	49033	4.806	ng	99
30) 1,2,4-Trichlorobenzene	10.799	180	61780	5.428	ng	98
31) Naphthalene	10.987	128	185958	5.385	ng	99
33) 4-Chloroaniline	11.105	127	66966	4.690	ng	98
34) Hexachlorobutadiene	11.269	225	38125	5.287	ng	96
36) 4-Chloro-3-methylphenol	12.216	107	51000	4.597	ng	94
37) 2-Methylnaphthalene	12.587	142	128538	5.237	ng	100
38) 1-Methylnaphthalene	12.805	142	119396	5.214	ng	99
40) 1,2,4,5-Tetrachloroben...	12.952	216	66909	5.285	ng	100
41) Hexachlorocyclopentadiene	12.928	237	26761	3.991	ng	92
43) 2,4,6-Trichlorophenol	13.193	196	38138	4.641	ng	93
44) 2,4,5-Trichlorophenol	13.263	196	45679	4.728	ng	95
46) 1,1'-Biphenyl	13.587	154	163817	5.359	ng	98
47) 2-Chloronaphthalene	13.634	162	123719	5.365	ng	98

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48) 2-Nitroaniline	13.840	65	28786	4.086	ng	95
49) Acenaphthylene	14.475	152	187495	5.081	ng	99
50) Dimethylphthalate	14.199	163	157209	5.208	ng	99
51) 2,6-Dinitrotoluene	14.322	165	29384	4.611	ng	98
52) Acenaphthene	14.816	154	117955	5.367	ng	96
53) 3-Nitroaniline	14.663	138	26010	3.993	ng	97
55) Dibenzofuran	15.146	168	194308	5.405	ng	99
57) 2,4-Dinitrotoluene	15.116	165	34526	4.120	ng	# 97
58) Fluorene	15.798	166	151793	5.466	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.375	232	37719	4.882	ng	99
60) Diethylphthalate	15.557	149	152773	5.081	ng	100
61) 4-Chlorophenyl-phenyle...	15.787	204	78073	5.491	ng	99
62) 4-Nitroaniline	15.822	138	24722m	5.637	ng	
63) Azobenzene	16.081	77	142380	4.963	ng	99
65) 4,6-Dinitro-2-methylph...	15.881	198	18340	3.787	ng	# 78
66) n-Nitrosodiphenylamine	16.004	169	126453	5.322	ng	99
67) 4-Bromophenyl-phenylether	16.687	248	45259	5.250	ng	99
68) Hexachlorobenzene	16.804	284	49786	5.244	ng	98
69) Atrazine	16.951	200	39124	5.157	ng	97
70) Pentachlorophenol	17.157	266	28257	4.343	ng	98
71) Phenanthrene	17.545	178	234110	5.509	ng	99
72) Anthracene	17.640	178	223059	5.241	ng	99
73) Carbazole	17.910	167	189187	5.020	ng	99
74) Di-n-butylphthalate	18.439	149	234084	4.837	ng	99
75) Fluoranthene	19.498	202	261745	5.273	ng	99
77) Benzidine	19.681	184	48261m	4.031	ng	
78) Pyrene	19.851	202	273295	5.073	ng	98
80) Butylbenzylphthalate	20.710	149	101974	4.467	ng	95
81) Benzo(a)anthracene	21.563	228	268773	5.138	ng	99
82) 3,3'-Dichlorobenzidine	21.492	252	75257	4.658	ng	97
83) Chrysene	21.622	228	270076	5.384	ng	99
84) Bis(2-ethylhexyl)phtha...	21.463	149	149399	4.443	ng	99
85) Di-n-octyl phthalate	22.439	149	252029	4.431	ng	98
87) Indeno(1,2,3-cd)pyrene	26.839	276	329000	5.292	ng	98
88) Benzo(b)fluoranthene	23.351	252	268634	5.151	ng	99
89) Benzo(k)fluoranthene	23.398	252	274415	5.225	ng	99
90) Benzo(a)pyrene	24.022	252	255497	5.074	ng	98
91) Dibenzo(a,h)anthracene	26.857	278	270446	5.297	ng	100
92) Benzo(g,h,i)perylene	27.668	276	270026	5.205	ng	98

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

