

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032122\
 Data File : BM034284.D
 Acq On : 21 Mar 2022 09:56
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 03/22/2022
 Supervised By : Jagrut Upadhyay 03/22/2022

Quant Time: Mar 21 17:53:43 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Mar 16 02:38:26 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	190920	20.000	ng	0.00
21) Naphthalene-d8	10.760	136	772950	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	501515	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	1037296	20.000	ng	0.00
76) Chrysene-d12	21.495	240	1086697	20.000	ng	0.00
86) Perylene-d12	23.912	264	1048665	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.490	112	865608	81.336	ng	-0.01
7) Phenol-d6	7.101	99	1204591	79.055	ng	0.00
23) Nitrobenzene-d5	9.107	82	1264742	79.664	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	544164	80.500	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	2899473	79.003	ng	0.00
79) Terphenyl-d14	19.942	244	4621345	73.895	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.349	88	194249	40.342	ng	99
3) Pyridine	3.755	79	522869	40.129	ng	98
4) n-Nitrosodimethylamine	3.666	42	245618	39.700	ng	99
6) Aniline	7.266	93	680914	39.104	ng	100
8) 2-Chlorophenol	7.507	128	494184	39.701	ng	99
9) Benzaldehyde	7.072	77	303263	36.941	ng	99
10) Phenol	7.125	94	597953	39.457	ng	99
11) bis(2-Chloroethyl)ether	7.360	93	481646	38.721	ng	100
12) 1,3-Dichlorobenzene	7.831	146	552390	39.122	ng	97
13) 1,4-Dichlorobenzene	7.978	146	560559	38.780	ng	98
14) 1,2-Dichlorobenzene	8.301	146	543840	38.594	ng	99
15) Benzyl Alcohol	8.184	79	475760	39.234	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.478	45	657784	38.534	ng	100
17) 2-Methylphenol	8.384	107	417697	39.354	ng	99
18) Hexachloroethane	9.037	117	209975	39.385	ng	98
19) n-Nitroso-di-n-propyla...	8.754	70	419410	38.053	ng	99
20) 3+4-Methylphenols	8.719	107	571913	38.962	ng	99
22) Acetophenone	8.766	105	772431	38.941	ng	# 97
24) Nitrobenzene	9.148	77	588791	39.723	ng	99
25) Isophorone	9.678	82	1034243	38.930	ng	98
26) 2-Nitrophenol	9.866	139	265808	40.438	ng	98
27) 2,4-Dimethylphenol	9.925	122	409507	39.786	ng	97
28) bis(2-Chloroethoxy)met...	10.160	93	669973	39.390	ng	99
29) 2,4-Dichlorophenol	10.401	162	482220	40.323	ng	98
30) 1,2,4-Trichlorobenzene	10.619	180	532177	39.149	ng	99
31) Naphthalene	10.807	128	1570917	39.011	ng	99
32) Benzoic acid	10.066	122	320336	37.636	ng	99
33) 4-Chloroaniline	10.913	127	638365	40.187	ng	99
34) Hexachlorobutadiene	11.101	225	338074	39.265	ng	99
35) Caprolactam	11.695	113	163683	39.374	ng	98
36) 4-Chloro-3-methylphenol	12.036	107	546345	39.866	ng	99
37) 2-Methylnaphthalene	12.419	142	1130116	39.330	ng	99
38) 1-Methylnaphthalene	12.636	142	1104511	39.295	ng	100
40) 1,2,4,5-Tetrachloroben...	12.789	216	586986	38.784	ng	99
41) Hexachlorocyclopentadiene	12.772	237	346970	39.594	ng	99
43) 2,4,6-Trichlorophenol	13.019	196	413315	39.556	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032122\
 Data File : BM034284.D
 Acq On : 21 Mar 2022 09:56
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 21 17:53:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 03/22/2022
 Supervised By :Jagrut Upadhyay 03/22/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.089	196	446441	39.806	ng	99
46) 1,1'-Biphenyl	13.425	154	1497511	39.161	ng	99
47) 2-Chloronaphthalene	13.466	162	1142876	38.937	ng	98
48) 2-Nitroaniline	13.660	65	366784	40.874	ng	99
49) Acenaphthylene	14.307	152	1824612	39.883	ng	100
50) Dimethylphthalate	14.042	163	1487134	38.960	ng	100
51) 2,6-Dinitrotoluene	14.154	165	318020	40.620	ng	99
52) Acenaphthene	14.648	154	1221872m	39.651	ng	
53) 3-Nitroaniline	14.483	138	325647	41.458	ng	98
54) 2,4-Dinitrophenol	14.683	184	183218	41.769	ng	97
55) Dibenzofuran	14.983	168	1804435	39.293	ng	98
56) 4-Nitrophenol	14.783	139	262101	41.160	ng	99
57) 2,4-Dinitrotoluene	14.942	165	435895	41.189	ng	94
58) Fluorene	15.630	166	1487421	39.897	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.207	232	397328	39.287	ng	99
60) Diethylphthalate	15.401	149	1561170	39.866	ng	100
61) 4-Chlorophenyl-phenyle...	15.624	204	749138	39.358	ng	100
62) 4-Nitroaniline	15.642	138	326071	43.153	ng	97
63) Azobenzene	15.919	77	1538783	40.749	ng	98
65) 4,6-Dinitro-2-methylph...	15.701	198	277675	41.168	ng	95
66) n-Nitrosodiphenylamine	15.836	169	1279607	38.829	ng	99
67) 4-Bromophenyl-phenylether	16.518	248	464684	38.969	ng	97
68) Hexachlorobenzene	16.636	284	517089	38.878	ng	99
69) Atrazine	16.783	200	451463	41.562	ng	100
70) Pentachlorophenol	16.977	266	340417	40.023	ng	98
71) Phenanthrene	17.365	178	2321469	39.963	ng	100
72) Anthracene	17.460	178	2281431	40.350	ng	100
73) Carbazole	17.724	167	2171332	41.177	ng	99
74) Di-n-butylphthalate	18.295	149	2729383	41.636	ng	99
75) Fluoranthene	19.377	202	2700034	41.462	ng	98
77) Benzidine	19.559	184	703174m	44.156	ng	
78) Pyrene	19.742	202	2824835	36.895	ng	100
80) Butylbenzylphthalate	20.636	149	1289650	39.829	ng	98
81) Benzo(a)anthracene	21.483	228	2696605	39.316	ng	100
82) 3,3'-Dichlorobenzidine	21.412	252	921043	40.205	ng	100
83) Chrysene	21.536	228	2697949	38.519	ng	100
84) Bis(2-ethylhexyl)phtha...	21.412	149	1924393	40.247	ng	98
85) Di-n-octyl phthalate	22.336	149	3191473	42.049	ng	100
87) Indeno(1,2,3-cd)pyrene	26.435	276	2543831	37.709	ng	98
88) Benzo(b)fluoranthene	23.177	252	2494045	39.772	ng	99
89) Benzo(k)fluoranthene	23.224	252	2685581	40.429	ng	100
90) Benzo(a)pyrene	23.806	252	2098549	39.476	ng	99
91) Dibenzo(a,h)anthracene	26.447	278	2092926	38.263	ng	99
92) Benzo(g,h,i)perylene	27.206	276	2079906	36.737	ng	99

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

