

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
 Data File : BM030329.D
 Acq On : 08 Jun 2021 18:48
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:54:04 PM

Quant Time: Jun 08 19:17:43 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 17:34:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	227932	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	864911	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	579514	20.000	ng	0.00
64) Phenanthrene-d10	17.209	188	1237937	20.000	ng	# 0.00
76) Chrysene-d12	21.374	240	1296808	20.000	ng	0.00
86) Perylene-d12	23.650	264	1106484	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	2334960	213.583	ng	0.00
7) Phenol-d6	7.016	99	3439659	231.718	ng	0.00
23) Nitrobenzene-d5	9.016	82	3193293	213.329	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0d	0.000	ng	
45) 2-Fluorobiphenyl	13.104	172	7612901	170.362	ng	0.00
79) Terphenyl-d14	19.827	244	11743182	166.206	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.352	88	471395	113.611	ng	90
3) Pyridine	3.746	79	1350651m	163.587	ng	
4) n-Nitrosodimethylamine	3.657	42	638021	154.779	ng	83
6) Aniline	7.187	93	1928711	120.575	ng	96
8) 2-Chlorophenol	7.416	128	1382895	102.967	ng	96
9) Benzaldehyde	6.993	77	583838	100.716	ng	94
10) Phenol	7.045	94	1711404	120.593	ng	93
11) bis(2-Chloroethyl)ether	7.281	93	1319875	118.548	ng	95
12) 1,3-Dichlorobenzene	7.745	146	1481880	90.995	ng	98
13) 1,4-Dichlorobenzene	7.887	146	1503095	90.369	ng	99
14) 1,2-Dichlorobenzene	8.204	146	1478677	90.659	ng	99
15) Benzyl Alcohol	8.092	79	1250109	120.478	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.381	45	2048688	169.872	ng	90
17) 2-Methylphenol	8.287	107	1139188	112.504	ng	93
18) Hexachloroethane	8.934	117	563893	97.057	ng	97
19) n-Nitroso-di-n-propyla...	8.669	70	1151970	120.252	ng	100
20) 3+4-Methylphenols	8.622	107	1555797	114.587	ng	97
22) Acetophenone	8.675	105	2020301	103.578	ng	# 98
24) Nitrobenzene	9.057	77	1629794	108.108	ng	96
25) Isophorone	9.581	82	3134426	113.333	ng	99
27) 2,4-Dimethylphenol	9.816	122	1190137	107.375	ng	96
28) bis(2-Chloroethoxy)met...	10.057	93	1684677	114.482	ng	100
29) 2,4-Dichlorophenol	10.292	162	1393418	91.176	ng	98
30) 1,2,4-Trichlorobenzene	10.510	180	1506333	81.128	ng	100
31) Naphthalene	10.692	128	4367924	98.708	ng	100
33) 4-Chloroaniline	10.804	127	1881105	111.275	ng	99
34) Hexachlorobutadiene	10.980	225	920091	70.375	ng	96
35) Caprolactam	11.616	113	447789m	120.727	ng	
36) 4-Chloro-3-methylphenol	11.922	107	1453215	109.981	ng	99
37) 2-Methylnaphthalene	12.304	142	3312882	97.378	ng	99
38) 1-Methylnaphthalene	12.522	142	3139202	96.867	ng	99
40) 1,2,4,5-Tetrachloroben...	12.669	216	1677476	79.297	ng	99
41) Hexachlorocyclopentadiene	12.651	237	972148	91.564	ng	98
43) 2,4,6-Trichlorophenol	12.910	196	1205014	89.855	ng	100
44) 2,4,5-Trichlorophenol	12.980	196	1306550	90.707	ng	99
46) 1,1'-Biphenyl	13.316	154	4074341	93.118	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
 Data File : BM030329.D
 Acq On : 08 Jun 2021 18:48
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:54:04 PM

Quant Time: Jun 08 19:17:43 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 17:34:01 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.357	162	3323743	94.150	ng	97
48) 2-Nitroaniline	13.557	65	993303	136.168	ng	99
49) Acenaphthylene	14.204	152	5236969	98.248	ng	99
50) Dimethylphthalate	13.939	163	4137258	93.483	ng	99
51) 2,6-Dinitrotoluene	14.057	165	971437	96.908	ng	95
52) Acenaphthene	14.545	154	3273077	99.081	ng	98
53) 3-Nitroaniline	14.386	138	985892	123.851	ng	99
54) 2,4-Dinitrophenol	14.586	184	588937	102.553	ng	88
55) Dibenzofuran	14.874	168	5158122	92.685	ng	99
58) Fluorene	15.521	166	4416073	96.624	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.098	232	1154432	86.228	ng	98
60) Diethylphthalate	15.298	149	4260048	98.397	ng	100
61) 4-Chlorophenyl-phenyle...	15.515	204	2204401	86.682	ng	95
62) 4-Nitroaniline	15.551	138	1052546	130.090	ng	88
63) Azobenzene	15.810	77	4182508	118.090	ng	98
65) 4,6-Dinitro-2-methylph...	15.604	198	830916	96.689	ng	93
66) n-Nitrosodiphenylamine	15.733	169	3826165	104.015	ng	99
67) 4-Bromophenyl-phenylether	16.404	248	1346870	91.388	ng	100
68) Hexachlorobenzene	16.527	284	1504520	91.455	ng	96
69) Atrazine	16.674	200	1011724	70.976	ng	96
71) Phenanthrene	17.257	178	6745625	96.692	ng	98
72) Anthracene	17.345	178	6747150	97.406	ng	98
73) Carbazole	17.615	167	6469952	103.566	ng	98
75) Fluoranthene	19.262	202	7959089	92.904	ng	94
77) Benzidine	19.439	184	2595489	108.205	ng	98
78) Pyrene	19.621	202	8246142	99.155	ng	96
80) Butylbenzylphthalate	20.509	149	3508571	110.625	ng	99
81) Benzo(a)anthracene	21.356	228	8008930	94.928	ng	94
82) 3,3'-Dichlorobenzidine	21.292	252	2726022	101.635	ng	100
83) Chrysene	21.409	228	7641051	91.264	ng	94
84) Bis(2-ethylhexyl)phtha...	21.280	149	5258262	104.030	ng	99
85) Di-n-octyl phthalate	22.168	149	8159089	96.115	ng	98
87) Indeno(1,2,3-cd)pyrene	25.985	276	7371620	88.010	ng	97
88) Benzo(b)fluoranthene	22.968	252	7642300	104.426	ng	97
89) Benzo(k)fluoranthene	23.015	252	7133945	100.610	ng	# 96
90) Benzo(a)pyrene	23.556	252	6702045	100.061	ng	# 98
91) Dibenzo(a,h)anthracene	25.997	278	6330257	87.546	ng	98
92) Benzo(g,h,i)perylene	26.697	276	5867436	85.974	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
 Data File : BM030329.D
 Acq On : 08 Jun 2021 18:48
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:54:04 PM

Quant Time: Jun 08 19:17:43 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 17:34:01 2021
 Response via : Initial Calibration

