

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052521\
 Data File : BF124048.D
 Acq On : 25 May 2021 16:19
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 5/26/2021 4:28:58 PM

Quant Time: May 25 17:10:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 14 19:05:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	51250	20.00	ng	0.00
21) Naphthalene-d8	8.181	136	183743	20.00	ng	# 0.00
39) Acenaphthene-d10	9.939	164	101225	20.00	ng	0.00
64) Phenanthrene-d10	11.427	188	179728	20.00	ng	0.00
76) Chrysene-d12	14.074	240	109711	20.00	ng	0.00
86) Perylene-d12	15.568	264	87991	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	283239	77.60	ng	0.00
7) Phenol-d6	6.528	99	365561	77.20	ng	0.00
23) Nitrobenzene-d5	7.463	82	362761	85.06	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	109150	89.78	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	631937	76.82	ng	0.00
79) Terphenyl-d14	13.016	244	630606	87.05	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.693	88	59566	37.16	ng	97
3) Pyridine	3.457	79	153369	37.07	ng	90
4) n-Nitrosodimethylamine	3.416	42	96028	35.73	ng	91
6) Aniline	6.563	93	202929	37.75	ng	96
8) 2-Chlorophenol	6.687	128	150680	39.67	ng	96
9) Benzaldehyde	6.445	77	86285	42.11	ng	95
10) Phenol	6.539	94	193701	40.47	ng	90
11) bis(2-Chloroethyl)ether	6.634	93	139757	37.57	ng	93
12) 1,3-Dichlorobenzene	6.839	146	176522	39.89	ng	95
13) 1,4-Dichlorobenzene	6.916	146	169608	37.46	ng	98
14) 1,2-Dichlorobenzene	7.069	146	173224	40.94	ng	95
15) Benzyl Alcohol	7.039	79	159012	38.88	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.175	45	260310	34.76	ng	98
17) 2-Methylphenol	7.145	107	125127	40.62	ng	94
18) Hexachloroethane	7.410	117	71666	42.71	ng	91
19) n-Nitroso-di-n-propyla...	7.310	70	124350	38.86	ng	93
20) 3+4-Methylphenols	7.298	107	155448	38.31	ng	90
22) Acetophenone	7.310	105	201526	38.62	ng	# 97
24) Nitrobenzene	7.481	77	180017	41.28	ng	97
25) Isophorone	7.722	82	318289	38.80	ng	97
26) 2-Nitrophenol	7.798	139	79884	48.70	ng	94
27) 2,4-Dimethylphenol	7.828	122	115069	36.64	ng	96
28) bis(2-Chloroethoxy)met...	7.928	93	161523	39.16	ng	96
29) 2,4-Dichlorophenol	8.039	162	131309	41.30	ng	89
30) 1,2,4-Trichlorobenzene	8.122	180	140977	39.44	ng	99
31) Naphthalene	8.204	128	426448	39.22	ng	99
32) Benzoic acid	7.951	122	53933m	31.29	ng	
33) 4-Chloroaniline	8.251	127	162193	39.18	ng	95
34) Hexachlorobutadiene	8.322	225	98041	41.90	ng	97
35) Caprolactam	8.622	113	35183	40.65	ng	# 83
36) 4-Chloro-3-methylphenol	8.728	107	138222	39.76	ng	# 86
37) 2-Methylnaphthalene	8.898	142	290908	39.47	ng	97
38) 1-Methylnaphthalene	8.992	142	272959	39.69	ng	94
40) 1,2,4,5-Tetrachloroben...	9.063	216	140196	38.40	ng	97
41) Hexachlorocyclopentadiene	9.045	237	82554	35.55	ng	92
43) 2,4,6-Trichlorophenol	9.169	196	92220	38.93	ng	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052521\
 Data File : BF124048.D
 Acq On : 25 May 2021 16:19
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 5/26/2021 4:28:58 PM

Quant Time: May 25 17:10:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 14 19:05:57 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	99026	40.24	ng	96
46) 1,1'-Biphenyl	9.357	154	338527	37.97	ng	97
47) 2-Chloronaphthalene	9.386	162	275382	38.89	ng	97
48) 2-Nitroaniline	9.475	65	105167	43.98	ng	92
49) Acenaphthylene	9.804	152	407727	38.97	ng	100
50) Dimethylphthalate	9.657	163	319003	39.47	ng	100
51) 2,6-Dinitrotoluene	9.722	165	71431	43.70	ng	87
52) Acenaphthene	9.975	154	285966	39.93	ng	95
53) 3-Nitroaniline	9.886	138	73774	45.41	ng	94
54) 2,4-Dinitrophenol	9.998	184	33471	41.91	ng	# 81
55) Dibenzofuran	10.145	168	408398	39.70	ng	97
56) 4-Nitrophenol	10.045	139	53578	39.87	ng	91
57) 2,4-Dinitrotoluene	10.127	165	97495	48.63	ng	93
58) Fluorene	10.486	166	308818	39.84	ng	91
59) 2,3,4,6-Tetrachlorophenol	10.263	232	76742	37.92	ng	# 97
60) Diethylphthalate	10.357	149	314170	39.25	ng	# 95
61) 4-Chlorophenyl-phenyle...	10.480	204	152504	39.26	ng	94
62) 4-Nitroaniline	10.504	138	70876	43.87	ng	91
63) Azobenzene	10.639	77	342845	38.51	ng	95
65) 4,6-Dinitro-2-methylph...	10.533	198	52964	46.00	ng	92
66) n-Nitrosodiphenylamine	10.598	169	271124	39.48	ng	97
67) 4-Bromophenyl-phenylether	10.969	248	95113	40.27	ng	99
68) Hexachlorobenzene	11.039	284	109468	41.51	ng	95
69) Atrazine	11.122	200	80796	41.66	ng	94
70) Pentachlorophenol	11.227	266	51776	33.12	ng	96
71) Phenanthrene	11.451	178	432118	37.94	ng	99
72) Anthracene	11.504	178	445163	38.57	ng	99
73) Carbazole	11.657	167	381321	37.76	ng	97
74) Di-n-butylphthalate	11.986	149	467343	36.98	ng	98
75) Fluoranthene	12.645	202	435746	36.94	ng	98
77) Benzidine	12.757	184	106421	40.20	ng	# 95
78) Pyrene	12.874	202	424586	43.59	ng	97
80) Butylbenzylphthalate	13.486	149	169052	44.26	ng	96
81) Benzo(a)anthracene	14.063	228	313185	39.47	ng	97
82) 3,3'-Dichlorobenzidine	14.021	252	102439	41.56	ng	99
83) Chrysene	14.104	228	302605	39.65	ng	97
84) Bis(2-ethylhexyl)phtha...	14.045	149	214722	39.84	ng	95
85) Di-n-octyl phthalate	14.662	149	303892	37.30	ng	98
87) Indeno(1,2,3-cd)pyrene	17.092	276	269168	39.61	ng	97
88) Benzo(b)fluoranthene	15.127	252	268993	39.52	ng	95
89) Benzo(k)fluoranthene	15.162	252	243524	39.26	ng	99
90) Benzo(a)pyrene	15.504	252	229950	39.08	ng	95
91) Dibenzo(a,h)anthracene	17.104	278	228411	39.91	ng	98
92) Benzo(g,h,i)perylene	17.551	276	226162	39.41	ng	97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052521\
Data File : BF124048.D
Acq On : 25 May 2021 16:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSTDCCC040

Manual Integrations
APPROVED

mohammad
5/26/2021 4:28:58 PM

Quant Time: May 25 17:10:30 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 14 19:05:57 2021
Response via : Initial Calibration

