

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061520\
 Data File : BP002411.D
 Acq On : 15 Jun 2020 16:59
 Operator : CG/JU
 Sample : L2182-08
 Misc : L00-WTR-10PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 LOQ-WATER-02-QT2-2020

Manual Integrations
 APPROVED

mohammad
 6/16/2020 11:07:27 AM

Quant Time: Jun 15 17:38:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 17:29:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	165364	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	661439	20.00	ng	0.00
39) Acenaphthene-d10	14.24	164	408992	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	870840	20.00	ng	0.00
76) Chrysene-d12	21.10	240	821162	20.00	ng	0.00
86) Perylene-d12	23.31	264	914966	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	1442693	161.44	ng	0.00
7) Phenol-d6	6.78	99	1972002	165.74	ng	0.00
23) Nitrobenzene-d5	8.74	82	1363026	123.06	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	682250	174.23	ng	0.00
45) 2-Fluorobiphenyl	12.86	172	2730139	112.24	ng	0.00
79) Terphenyl-d14	19.59	244	3708798	118.34	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.16	88	44443	10.796	ng	99
3) Pyridine	3.55	79	106608	9.528	ng	98
4) n-Nitrosodimethylamine	3.46	42	46359	10.055	ng	100
6) Aniline	6.92	93	172713	10.674	ng	99
8) 2-Chlorophenol	7.16	128	102419	10.193	ng	97
9) Benzaldehyde	6.73	77	100421	16.519	ng	96
10) Phenol	6.80	94	137328	10.642	ng	99
11) bis(2-Chloroethyl)ether	7.03	93	112573	10.449	ng	99
12) 1,3-Dichlorobenzene	7.48	146	117254	10.134	ng	99
13) 1,4-Dichlorobenzene	7.63	146	119393	10.256	ng	99
14) 1,2-Dichlorobenzene	7.94	146	113210	10.152	ng	100
15) Benzyl Alcohol	7.83	79	87589	9.498	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.13	45	156545	10.240	ng	99
17) 2-Methylphenol	8.04	107	87973	9.330	ng	99
18) Hexachloroethane	8.66	117	43420	10.238	ng	93
19) n-Nitroso-di-n-propylamine	8.39	70	85908	10.802	ng	97
20) 3+4-Methylphenols	8.37	107	118999	9.351	ng	98
22) Acetophenone	8.40	105	172526	11.610	ng	99
24) Nitrobenzene	8.78	77	136313	11.449	ng	95
25) Isophorone	9.30	82	239356	10.611	ng	99
26) 2-Nitrophenol	9.49	139	51788	9.761	ng	97
27) 2,4-Dimethylphenol	9.56	122	78224	9.399	ng	99
28) bis(2-Chloroethoxy)methane	9.79	93	153080	11.393	ng	98
29) 2,4-Dichlorophenol	10.02	162	93796	10.011	ng	97
30) 1,2,4-Trichlorobenzene	10.23	180	107841	10.326	ng	99
31) Naphthalene	10.42	128	344068	11.197	ng	99
32) Benzoic acid	9.65	122	28654m	4.525	ng	
33) 4-Chloroaniline	10.53	127	140032	10.438	ng	99
34) Hexachlorobutadiene	10.72	225	60249	9.886	ng	94
35) Caprolactam	11.26	113	34135	10.330	ng	94
36) 4-Chloro-3-methylphenol	11.67	107	102319	10.093	ng	97
37) 2-Methylnaphthalene	12.04	142	245953	11.277	ng	99
38) 1-Methylnaphthalene	12.26	142	232196	11.342	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.42	216	124621	11.559	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061520\
 Data File : BP002411.D
 Acq On : 15 Jun 2020 16:59
 Operator : CG/JU
 Sample : L2182-08
 Misc : L00-WTR-10PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 LOQ-WATER-02-QT2-2020

Manual Integrations
 APPROVED

mohammad
 6/16/2020 11:07:27 AM

Quant Time: Jun 15 17:38:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 17:29:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.40	237	47188	8.233	nq	98
43) 2,4,6-Trichlorophenol	12.66	196	73484	9.631	nq	95
44) 2,4,5-Trichlorophenol	12.73	196	80016	9.056	nq	99
46) 1,1'-Biphenyl	13.07	154	326740	12.190	nq	99
47) 2-Chloronaphthalene	13.10	162	232565	10.607	nq	98
48) 2-Nitroaniline	13.31	65	64819	9.261	nq	96
49) Acenaphthylene	13.96	152	381653	10.906	nq	98
50) Dimethylphthalate	13.70	163	295547	10.546	nq	99
51) 2,6-Dinitrotoluene	13.82	165	62020	10.191	nq	99
52) Acenaphthene	14.30	154	226821	11.064	nq	98
53) 3-Nitroaniline	14.15	138	61098	8.792	nq	98
54) 2,4-Dinitrophenol	14.36	184	20103	6.022	nq	97
55) Dibenzofuran	14.65	168	372467	11.285	nq	98
56) 4-Nitrophenol	14.47	139	46125	8.359	nq	98
57) 2,4-Dinitrotoluene	14.61	165	78288	9.450	nq	95
58) Fluorene	15.30	166	294724	11.352	nq	95
59) 2,3,4,6-Tetrachlorophenol	14.87	232	70873	10.105	nq	99
60) Diethylphthalate	15.09	149	293476	10.452	nq	98
61) 4-Chlorophenyl-phenylether	15.30	204	145753	11.582	nq	98
62) 4-Nitroaniline	15.31	138	61912	9.272	nq	96
63) Azobenzene	15.59	77	319355	11.426	nq	95
65) 4,6-Dinitro-2-methylphenol	15.38	198	39082	8.505	nq	96
66) n-Nitrosodiphenylamine	15.52	169	260797	11.723	nq	97
67) 4-Bromophenyl-phenylether	16.20	248	83095	10.073	nq	97
68) Hexachlorobenzene	16.32	284	94775	10.827	nq	96
69) Atrazine	16.47	200	90428	12.461	nq	100
70) Pentachlorophenol	16.66	266	44730	8.541	nq	97
71) Phenanthrene	17.04	178	473312	11.619	nq	98
72) Anthracene	17.13	178	464714	11.484	nq	98
73) Carbazole	17.40	167	428938	10.775	nq	100
74) Di-n-butylphthalate	17.99	149	484257	10.279	nq	97
75) Fluoranthene	19.03	202	538840	11.082	nq	99
77) Benzidine	19.21	184	257564	14.037	nq	97
78) Pyrene	19.37	202	581319	11.557	nq	98
80) Butylbenzylphthalate	20.27	149	199256	8.912	nq	99
81) Benzo(a)anthracene	21.09	228	541183	11.546	nq	97
82) 3,3'-Dichlorobenzidine	21.03	252	202451	11.622	nq	99
83) Chrysene	21.14	228	530495	11.946	nq	98
84) Bis(2-ethylhexyl)phthalate	21.04	149	327540	10.071	nq	99
85) Di-n-octyl phthalate	21.92	149	536900	9.412	nq	99
87) Indeno(1,2,3-cd)pyrene	25.52	276	606304	10.594	nq	98
88) Benzo(b)fluoranthene	22.64	252	530820	10.758	nq	97
89) Benzo(k)fluoranthene	22.69	252	535431	11.421	nq	99
90) Benzo(a)pyrene	23.21	252	479176	10.363	nq	98
91) Dibenzo(a,h)anthracene	25.53	278	509124	11.091	nq	96
92) Benzo(g,h,i)perylene	26.19	276	499942	10.515	nq	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061520\
Data File : BP002411.D
Acq On : 15 Jun 2020 16:59
Operator : CG/JU
Sample : L2182-08
Misc : LOO-WTR-10PPM
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
LOQ-WATER-02-QT2-2020

Manual Integrations
APPROVED

mohammad
6/16/2020 11:07:27 AM

Quant Time: Jun 15 17:38:06 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 15 17:29:37 2020
Response via : Initial Calibration

