Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006297.D
Acq On : 23 Jul 2021 21:00
Operator : CG/JU
Sample : SSTD16006
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SSTD160606

Manual Integrations APPROVED
mohammad
7/26/2021 1:29:46 PM

Quant Time: Jul 24 02:47:25 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 24 02:41:40 2021
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	306404	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1357495	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	798091	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1587681	20.000	ng/ul	0.01
79) Chrysene-d12	21.280	240	1485230	20.000	ng/ul	0.01
88) Perylene-d12	23.633	264	1541623	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.705	84	4116808	217.652	ng/ul	0.00
7) Phenol-d5	6.981	99	4873088	195.861	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.146	67	2858402	217.881	ng/ul	0.01
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/ul	
15) 4-Methylphenol-d8	8.522	113	3756656	190.364	ng/ul	0.02
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	10.734	131	5067923	165.140	ng/ul	0.02
46) Dimethylphthalate-d6	0.000	166	0d	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0d	0.000	ng/ul	
54) 4-Nitrophenol-d4	14.663	143	2040798	190.337	ng/ul	0.05
60) Fluorene-d10	0.000	176	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	15.563	200	1724697	155.605	ng/ul	0.03
73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	
Target Compounds						Qvalue
5) Pyridine	3.723	79	4227985	216.625	ng/ul	99
6) Benzaldehyde	6.946	77	2062227	159.442	ng/ul	100
8) Phenol	7.011	94	4846127	186.684	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.240	93	3685910	186.842	ng/ul	99
13) 2-Methylphenol	8.246	108	3580421	181.130	ng/ul	100
14) 2,2'-oxybis(1-Chloropr...	8.340	45	4052133	206.706	ng/ul	97
16) Acetophenone	8.634	105	5480792	169.022	ng/ul	99
18) 4-Methylphenol	8.599	108	3813149	181.152	ng/ul	97
32) 4-Chloroaniline	10.757	127	4925407	158.655	ng/ul	99
34) Caprolactam	11.599	113	1202385m	163.972	ng/ul	
40) Hexachlorocyclopentadiene	12.593	237	2445378	134.277	ng/ul	98
51) 3-Nitroaniline	14.351	138	2061895	174.703	ng/ul	97
53) 2,4-Dinitrophenol	14.551	184	1284383	154.196	ng/ul	93
55) 4-Nitrophenol	14.675	109	1565747	155.050	ng/ul	98
63) 4-Nitroaniline	15.534	138	1949533m	168.134	ng/ul	
66) 4,6-Dinitro-2-methylph...	15.575	198	1651017	148.064	ng/ul#	95
70) Atrazine	16.669	200	2636920	136.193	ng/ul	99
71) Pentachlorophenol	16.828	266	2011201	146.404	ng/ul	98
77) Carbazole	17.592	167	12111350	151.460	ng/ul	95
84) 3,3'-Dichlorobenzidine	21.204	252	5529308	162.002	ng/ul	97
89) Di-n-octyl phthalate	22.116	149	14603179	147.684	ng/ul	100

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

Instrument :
BNA_P
ClientSampleId :
SSTD160606

Manual Integrations APPROVED

mohammad
7/26/2021 1:29:46 PM

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006297.D
Acq On : 23 Jul 2021 21:00
Operator : CG/JU
Sample : SSTD16006
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Jul 24 02:47:25 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 24 02:41:40 2021
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
SSTD160606

Manual Integrations APPROVED
mohammad
7/26/2021 1:29:46 PM