

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012521\
 Data File : BN013479.D
 Acq On : 26 Jan 2021 00:16
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4007

Quant Time: Jan 26 00:45:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN012521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 25 22:48:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.574	152	4683	0.40	ng/ul	0.00
4) Naphthalene-d8	10.334	136	17031	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.204	164	8763	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	18191	0.40	ng/ul	0.00
17) Chrysene-d12	21.151	240	15659	0.40	ng/ul	0.00
23) Perylene-d12	23.320	264	14062	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.174	96	1856	0.39	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.929	152	9700	0.39	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	17899	0.38	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.207	88	1792	0.38	ng/ul	98
5) Naphthalene	10.378	128	18751	0.39	ng/ul	100
7) 2-Methylnaphthalene	12.001	142	12492	0.39	ng/ul	100
8) 1-Methylnaphthalene	12.226	142	11975	0.39	ng/ul	100
10) Acenaphthylene	13.922	152	14315	0.39	ng/ul	100
11) Acenaphthene	14.264	153	12237	0.39	ng/ul	99
12) Fluorene	15.259	166	13555	0.39	ng/ul	98
14) Pentachlorophenol	16.603	266	840	0.38	ng/ul	97
15) Phenanthrene	16.992	178	22495	0.38	ng/ul	100
16) Anthracene	17.085	178	18357	0.39	ng/ul	99
19) Fluoranthene	19.019	202	23717	0.38	ng/ul	99
20) Pyrene	19.381	202	23931	0.38	ng/ul	98
21) Benzo(a)anthracene	21.137	228	19138	0.39	ng/ul	99
22) Chrysene	21.186	228	23593	0.39	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	22346	0.36	ng/ul	99
25) Benzo(k)fluoranthene	22.717	252	23891	0.37	ng/ul	100
26) Benzo(a)pyrene	23.226	252	18130	0.38	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.480	276	24916	0.37	ng/ul#	81
28) Dibenzo(a,h)anthracene	25.493	278	20166	0.37	ng/ul	99
29) Benzo(g,h,i)perylene	26.134	276	21492	0.36	ng/ul	99

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

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