

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012621\
 Data File : BN013483.D
 Acq On : 26 Jan 2021 09:15
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4008

Quant Time: Jan 26 13:48:24 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 : Tue Jan 26 10:37:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.570	152	4708	0.40	ng/ul	0.00
4) Naphthalene-d8	10.329	136	17255	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.204	164	8869	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	18545	0.40	ng/ul	0.00
17) Chrysene-d12	21.154	240	15654	0.40	ng/ul	0.00
23) Perylene-d12	23.325	264	13654	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.174	96	1855	0.39	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.924	152	9829	0.39	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	18175	0.39	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.207	88	1799	0.37	ng/ul	96
5) Naphthalene	10.378	128	18877	0.39	ng/ul	99
7) 2-Methylnaphthalene	12.001	142	12683	0.39	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	12046	0.39	ng/ul	99
10) Acenaphthylene	13.917	152	14347	0.38	ng/ul	100
11) Acenaphthene	14.264	153	12421	0.39	ng/ul	100
12) Fluorene	15.254	166	13831	0.39	ng/ul	97
14) Pentachlorophenol	16.603	266	927	0.41	ng/ul	99
15) Phenanthrene	16.992	178	23095	0.39	ng/ul	100
16) Anthracene	17.084	178	18668	0.39	ng/ul	100
19) Fluoranthene	19.018	202	24327	0.39	ng/ul	100
20) Pyrene	19.381	202	24373	0.39	ng/ul	99
21) Benzo(a)anthracene	21.139	228	18973	0.38	ng/ul	99
22) Chrysene	21.189	228	23814	0.39	ng/ul	100
24) Benzo(b)fluoranthene	22.679	252	22329	0.37	ng/ul	99
25) Benzo(k)fluoranthene	22.723	252	23551	0.38	ng/ul	99
26) Benzo(a)pyrene	23.232	252	17724	0.38	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.480	276	24707	0.38	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.497	278	20139	0.38	ng/ul	100
29) Benzo(g,h,i)perylene	26.134	276	21884	0.38	ng/ul	97

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

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