

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012321\
 Data File : BN013467.D
 Acq On : 22 Jan 2021 22:22
 Operator : CG/JU
 Sample : SST0.806
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8006

Quant Time: Jan 23 01:16:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN012321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 23 01:14:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.574	152	3991	0.40	ng/ul	0.00
4) Naphthalene-d8	10.333	136	14755	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.207	164	7740	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.957	188	16394	0.40	ng/ul	0.00
17) Chrysene-d12	21.156	240	16469	0.40	ng/ul	0.00
23) Perylene-d12	23.330	264	13889	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	3377	0.75	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.933	152	17534	0.78	ng/ul	0.00
18) Fluoranthene-d10	18.990	212	35140	0.72	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.207	88	3412	0.68	ng/ul	91
5) Naphthalene	10.382	128	33840	0.81	ng/ul	99
7) 2-Methylnaphthalene	12.005	142	21986	0.80	ng/ul	100
8) 1-Methylnaphthalene	12.230	142	21132	0.81	ng/ul	100
10) Acenaphthylene	13.920	152	26305	0.84	ng/ul	100
11) Acenaphthene	14.267	153	22800	0.83	ng/ul	99
12) Fluorene	15.262	166	25694	0.82	ng/ul	98
14) Pentachlorophenol	16.611	266	1958	0.47	ng/ul	98
15) Phenanthrene	16.995	178	43009	0.85	ng/ul	99
16) Anthracene	17.088	178	36349	0.82	ng/ul	100
19) Fluoranthene	19.022	202	49057	0.76	ng/ul	100
20) Pyrene	19.385	202	48771	0.75	ng/ul	99
21) Benzo(a)anthracene	21.141	228	46468	0.86	ng/ul	99
22) Chrysene	21.191	228	49094	0.84	ng/ul	99
24) Benzo(b)fluoranthene	22.684	252	45844	0.82	ng/ul	98
25) Benzo(k)fluoranthene	22.725	252	48609	0.91	ng/ul	98
26) Benzo(a)pyrene	23.234	252	38369	0.83	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.492	276	52477	0.94	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.509	278	42893	0.91	ng/ul	98
29) Benzo(g,h,i)perylene	26.146	276	44131	0.90	ng/ul	99

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

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