

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012321\
 Data File : BN013466.D
 Acq On : 22 Jan 2021 21:47
 Operator : CG/JU
 Sample : SST0.405
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4005

Quant Time: Jan 23 01:16:26 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	4329	0.40	ng/ul	0.00
4) Naphthalene-d8	10.334	136	15820	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.209	164	8296	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.958	188	17641	0.40	ng/ul	0.00
17) Chrysene-d12	21.157	240	17425	0.40	ng/ul	0.00
23) Perylene-d12	23.328	264	14458	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	1805	0.37	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.929	152	8890	0.37	ng/ul	0.00
18) Fluoranthene-d10	18.990	212	17762	0.34	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.207	88	1711	0.31	ng/ul	100
5) Naphthalene	10.384	128	17525	0.39	ng/ul	100
7) 2-Methylnaphthalene	12.006	142	11271	0.38	ng/ul	100
8) 1-Methylnaphthalene	12.226	142	10807	0.38	ng/ul	100
10) Acenaphthylene	13.922	152	13338	0.40	ng/ul	100
11) Acenaphthene	14.269	153	11547	0.39	ng/ul	100
12) Fluorene	15.263	166	13020	0.39	ng/ul	100
14) Pentachlorophenol	16.612	266	961	0.21	ng/ul	100
15) Phenanthrene	16.996	178	22182	0.41	ng/ul	100
16) Anthracene	17.089	178	18410	0.39	ng/ul	100
19) Fluoranthene	19.023	202	24759	0.36	ng/ul	100
20) Pyrene	19.386	202	25246	0.37	ng/ul	100
21) Benzo(a)anthracene	21.139	228	23604	0.41	ng/ul	100
22) Chrysene	21.192	228	25284	0.41	ng/ul	100
24) Benzo(b)fluoranthene	22.682	252	22844	0.39	ng/ul	100
25) Benzo(k)fluoranthene	22.723	252	24671	0.44	ng/ul	100
26) Benzo(a)pyrene	23.235	252	19561	0.41	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.490	276	25452	0.44	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.503	278	20756	0.42	ng/ul	100
29) Benzo(g,h,i)perylene	26.147	276	22087	0.43	ng/ul	100

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

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