

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061721\
 Data File : BN014936.D
 Acq On : 17 Jun 2021 10:28
 Operator : CG/JU
 Sample : SST0.806
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8006

Quant Time: Jun 17 11:04:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 17 11:01:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.747	152	6101	0.400	ng/ul	0.00
4) Naphthalene-d8	10.516	136	23008	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.371	164	13142	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.118	188	27256	0.400	ng/ul	0.00
17) Chrysene-d12	21.318	240	25244	0.400	ng/ul	0.00
23) Perylene-d12	23.574	264	21724	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	7010	1.120	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.111	152	30970	0.778	ng/ul	0.00
18) Fluoranthene-d10	19.153	212	65580	1.030	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	6573	1.005	ng/ul#	93
5) Naphthalene	10.565	128	57405	0.859	ng/ul	100
7) 2-Methylnaphthalene	12.182	142	37834	0.787	ng/ul	100
8) 1-Methylnaphthalene	12.402	142	37092	0.799	ng/ul	99
10) Acenaphthylene	14.088	152	54725	1.151	ng/ul	100
11) Acenaphthene	14.431	153	38809	0.857	ng/ul	99
12) Fluorene	15.421	166	44697	0.804	ng/ul	99
14) Pentachlorophenol	16.764	266	5149	1.342	ng/ul	99
15) Phenanthrene	17.156	178	71468	0.843	ng/ul	99
16) Anthracene	17.249	178	68610	1.038	ng/ul	99
19) Fluoranthene	19.181	202	81054	1.033	ng/ul	99
20) Pyrene	19.544	202	81985	1.032	ng/ul	99
21) Benzo(a)anthracene	21.303	228	73248	0.918	ng/ul	100
22) Chrysene	21.356	228	78501	0.861	ng/ul	99
24) Benzo(b)fluoranthene	22.902	252	67643	0.708	ng/ul	99
25) Benzo(k)fluoranthene	22.945	252	79925	0.794	ng/ul	100
26) Benzo(a)pyrene	23.477	252	62165	0.852	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.856	276	74907	0.721	ng/ul	99
28) Dibenzo(a,h)anthracene	25.876	278	59885	0.719	ng/ul	100
29) Benzo(g,h,i)perylene	26.543	276	67690	0.764	ng/ul	100

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

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