

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062121\
 Data File : BN014951.D
 Acq On : 21 Jun 2021 12:13
 Operator : CG/JU
 Sample : SST0.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4006

Quant Time: Jun 21 13:16:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 17 13:23:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.747	152	3169	0.400	ng/ul	0.00
4) Naphthalene-d8	10.516	136	10721	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.371	164	5435	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.114	188	10902	0.400	ng/ul	0.00
17) Chrysene-d12	21.318	240	12365	0.400	ng/ul	0.00
23) Perylene-d12	23.574	264	9096	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	1445	0.321	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.110	152	6819	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.149	212	12115	0.311	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.333	88	1534	0.367	ng/ul#	81
5) Naphthalene	10.565	128	14403	0.424	ng/ul	99
7) 2-Methylnaphthalene	12.182	142	9404	0.421	ng/ul	99
8) 1-Methylnaphthalene	12.402	142	8705	0.403	ng/ul	100
10) Acenaphthylene	14.088	152	11704	0.409	ng/ul	99
11) Acenaphthene	14.431	153	9011	0.444	ng/ul	99
12) Fluorene	15.421	166	10040	0.432	ng/ul	99
14) Pentachlorophenol	16.764	266	1062	0.389	ng/ul	99
15) Phenanthrene	17.156	178	16500	0.454	ng/ul	99
16) Anthracene	17.249	178	15563	0.440	ng/ul	99
19) Fluoranthene	19.181	202	19733	0.398	ng/ul	95
20) Pyrene	19.544	202	20209	0.398	ng/ul	94
21) Benzo(a)anthracene	21.303	228	16907	0.378	ng/ul	99
22) Chrysene	21.353	228	22082	0.453	ng/ul	98
24) Benzo(b)fluoranthene	22.899	252	16865	0.473	ng/ul	100
25) Benzo(k)fluoranthene	22.945	252	19302	0.466	ng/ul	100
26) Benzo(a)pyrene	23.477	252	14081	0.431	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.852	276	17277	0.459	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.869	278	13721	0.458	ng/ul	96
29) Benzo(g,h,i)perylene	26.536	276	16038	0.464	ng/ul	98

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

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