

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080323\
 Data File : BN026802.D
 Acq On : 03 Aug 2023 12:09
 Operator : MA/JU
 Sample : SST0.220
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2238

Quant Time: Aug 04 03:38:16 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 03 15:46:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.118	152	8868	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	23641	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.744	164	9755	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.494	188	30473	0.400	ng/ul	0.00
17) Chrysene-d12	21.668	240	30552	0.400	ng/ul	0.00
23) Perylene-d12	24.223	264	35072	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.448	96	2112	0.201	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.517	152	6833	0.196	ng/ul	0.00
18) Fluoranthene-d10	19.506	212	19042	0.197	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	2074	0.193	ng/ul	91
5) Naphthalene	10.999	128	13200	0.198	ng/ul	99
7) 2-Methylnaphthalene	12.594	142	7972	0.197	ng/ul	100
8) 1-Methylnaphthalene	12.808	142	8359	0.197	ng/ul	100
10) Acenaphthylene	14.476	152	11127	0.204	ng/ul	99
11) Acenaphthene	14.809	153	7697	0.204	ng/ul	99
12) Fluorene	15.790	166	8369	0.196	ng/ul	99
14) Pentachlorophenol	17.114	266	1661	0.193	ng/ul	100
15) Phenanthrene	17.536	178	20291	0.203	ng/ul	100
16) Anthracene	17.624	178	17860	0.197	ng/ul	99
19) Fluoranthene	19.538	202	25088	0.198	ng/ul	99
20) Pyrene	19.901	202	27171	0.201	ng/ul	99
21) Benzo(a)anthracene	21.650	228	24470	0.201	ng/ul	98
22) Chrysene	21.708	228	24181	0.202	ng/ul	99
24) Benzo(b)fluoranthene	23.430	252	26425	0.199	ng/ul	96
25) Benzo(k)fluoranthene	23.483	252	26256	0.201	ng/ul	96
26) Benzo(a)pyrene	24.109	252	24199	0.202	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.928	276	31228	0.199	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.965	278	24936	0.200	ng/ul	95
29) Benzo(g,h,i)perylene	27.770	276	25998	0.199	ng/ul	97

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

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