

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082423\
 Data File : BM041459.D
 Acq On : 24 Aug 2023 10:22
 Operator : MA/JU
 Sample : SST0.207
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.2044

Quant Time: Aug 24 18:22:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 24 13:21:17 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.118	152	8659	0.400	ng/ul	0.00
4) Naphthalene-d8	10.950	136	23147	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.759	164	10378	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.515	188	20202	0.400	ng/ul	0.00
17) Chrysene-d12	21.705	240	8910	0.400	ng/ul	0.00
23) Perylene-d12	24.281	264	9636	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.452	96	2278	0.208	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.528	152	6209	0.203	ng/ul	0.00
18) Fluoranthene-d10	19.534	212	7618	0.194	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.490	88	2311	0.205	ng/ul#	85
5) Naphthalene	11.005	128	14251	0.207	ng/ul	99
7) 2-Methylnaphthalene	12.599	142	8398	0.203	ng/ul	99
8) 1-Methylnaphthalene	12.819	142	8529	0.202	ng/ul	99
10) Acenaphthylene	14.486	152	12875	0.207	ng/ul	98
11) Acenaphthene	14.823	153	8845	0.204	ng/ul	98
12) Fluorene	15.809	166	9791	0.202	ng/ul	100
14) Pentachlorophenol	17.130	266	775	0.203	ng/ul	97
15) Phenanthrene	17.557	178	15189	0.205	ng/ul	99
16) Anthracene	17.650	178	13676	0.203	ng/ul	98
19) Fluoranthene	19.566	202	15520	0.190	ng/ul	100
20) Pyrene	19.933	202	16036	0.194	ng/ul	100
21) Benzo(a)anthracene	21.688	228	11746	0.203	ng/ul	99
22) Chrysene	21.743	228	12755	0.204	ng/ul	100
24) Benzo(b)fluoranthene	23.483	252	13002	0.197	ng/ul	93
25) Benzo(k)fluoranthene	23.535	252	12926	0.196	ng/ul	97
26) Benzo(a)pyrene	24.164	252	10650	0.196	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	27.005	276	13575	0.200	ng/ul#	99
28) Dibenzo(a,h)anthracene	27.045	278	10443	0.201	ng/ul	95
29) Benzo(g,h,i)perylene	27.850	276	12050	0.202	ng/ul	97

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

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