

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062024\
 Data File : BM046175.D
 Acq On : 22 Jun 2024 01:02
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4139

Quant Time: Jun 22 01:31:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 18 13:58:29 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.480	152	1790	0.400	ng/ul	0.02
4) Naphthalene-d8	10.255	136	6178	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.100	164	3562	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.858	188	6417	0.400	ng/ul	0.00
17) Chrysene-d12	21.035	240	5231	0.400	ng/ul	0.00
23) Perylene-d12	23.124	264	7225	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.042	96	1025	0.417	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.871	152	3279	0.395	ng/ul	0.00
18) Fluoranthene-d10	18.881	212	6242	0.398	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.067	88	1142	0.415	ng/ul	96
5) Naphthalene	10.304	128	6494	0.377	ng/ul	97
7) 2-Methylnaphthalene	11.948	142	3683	0.380	ng/ul	98
8) 1-Methylnaphthalene	12.146	142	4351	0.409	ng/ul	98
10) Acenaphthylene	13.827	152	6403	0.379	ng/ul	99
11) Acenaphthene	14.164	153	4702	0.386	ng/ul	98
12) Fluorene	15.182	166	4872	0.372	ng/ul	99
14) Pentachlorophenol	16.546	266	705	0.367	ng/ul	100
15) Phenanthrene	16.901	178	6874	0.381	ng/ul	100
16) Anthracene	17.006	178	4414	0.329	ng/ul	98
19) Fluoranthene	18.914	202	8918	0.394	ng/ul	99
20) Pyrene	19.272	202	9687	0.395	ng/ul	97
21) Benzo(a)anthracene	21.020	228	7228	0.358	ng/ul	99
22) Chrysene	21.070	228	9979	0.387	ng/ul	99
24) Benzo(b)fluoranthene	22.507	252	9651	0.371	ng/ul	99
25) Benzo(k)fluoranthene	22.548	252	11990	0.387	ng/ul	99
26) Benzo(a)pyrene	23.039	252	8876	0.368	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.182	276	14307	0.362	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.195	278	10822	0.368	ng/ul	99
29) Benzo(g,h,i)perylene	25.802	276	12046	0.356	ng/ul	97

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

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