

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070323\
 Data File : BM040600.D
 Acq On : 04 Jul 2023 07:15
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4149

Quant Time: Jul 04 07:44:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 13:40:38 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.919	152	5483	0.400	ng/ul	0.00
4) Naphthalene-d8	10.722	136	20766	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.562	164	9651	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.314	188	16969	0.400	ng/ul	0.00
17) Chrysene-d12	21.499	240	9987	0.400	ng/ul	0.00
23) Perylene-d12	23.904	264	10552	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	2875	0.334	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.316	152	11290	0.388	ng/ul	0.00
18) Fluoranthene-d10	19.341	212	13772	0.430	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.350	88	3023	0.345	ng/ul#	82
5) Naphthalene	10.777	128	21760	0.380	ng/ul	99
7) 2-Methylnaphthalene	12.388	142	12904	0.384	ng/ul	98
8) 1-Methylnaphthalene	12.608	142	12445	0.370	ng/ul	99
10) Acenaphthylene	14.284	152	17042	0.395	ng/ul	99
11) Acenaphthene	14.622	153	13551	0.372	ng/ul	98
12) Fluorene	15.612	166	14190	0.365	ng/ul	100
14) Pentachlorophenol	16.968	266	773	0.237	ng/ul	90
15) Phenanthrene	17.356	178	19173	0.364	ng/ul	99
16) Anthracene	17.449	178	16794	0.379	ng/ul	99
19) Fluoranthene	19.374	202	17574	0.414	ng/ul	98
20) Pyrene	19.736	202	18145	0.399	ng/ul	99
21) Benzo(a)anthracene	21.481	228	13734	0.407	ng/ul	99
22) Chrysene	21.534	228	14141	0.367	ng/ul	99
24) Benzo(b)fluoranthene	23.168	252	14767	0.362	ng/ul	93
25) Benzo(k)fluoranthene	23.217	252	13948	0.347	ng/ul	92
26) Benzo(a)pyrene	23.799	252	12920	0.351	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.402	276	17674	0.345	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.419	278	14064	0.351	ng/ul	98
29) Benzo(g,h,i)perylene	27.170	276	15158	0.338	ng/ul	100

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

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