

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040905.D
 Acq On : 15 Jul 2023 08:39
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4180

Quant Time: Jul 16 00:28:37 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 : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.861	152	3506	0.400	ng/ul	0.00
4) Naphthalene-d8	10.658	136	11262	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.504	164	4573	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	7508	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	5320	0.400	ng/ul	0.00
23) Perylene-d12	23.827	264	5508	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	2331	0.423	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.253	152	5962	0.378	ng/ul	0.00
18) Fluoranthene-d10	19.287	212	6761	0.396	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	2382	0.425	ng/ul	92
5) Naphthalene	10.708	128	12393	0.399	ng/ul	99
7) 2-Methylnaphthalene	12.324	142	6791	0.372	ng/ul	98
8) 1-Methylnaphthalene	12.544	142	6796	0.373	ng/ul	99
10) Acenaphthylene	14.226	152	7709	0.377	ng/ul	99
11) Acenaphthene	14.564	153	6653	0.386	ng/ul	98
12) Fluorene	15.559	166	6489	0.352	ng/ul	100
14) Pentachlorophenol	16.919	266	610	0.422	ng/ul	91
15) Phenanthrene	17.299	178	8801	0.377	ng/ul	99
16) Anthracene	17.392	178	7117	0.363	ng/ul	98
19) Fluoranthene	19.320	202	8620	0.381	ng/ul	98
20) Pyrene	19.682	202	9012	0.372	ng/ul	99
21) Benzo(a)anthracene	21.434	228	6831	0.380	ng/ul	99
22) Chrysene	21.486	228	7808	0.380	ng/ul	99
24) Benzo(b)fluoranthene	23.102	252	8089	0.379	ng/ul	98
25) Benzo(k)fluoranthene	23.149	252	8048	0.383	ng/ul	96
26) Benzo(a)pyrene	23.719	252	7200	0.375	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.283	276	10570	0.395	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.297	278	8432	0.403	ng/ul	98
29) Benzo(g,h,i)perylene	27.031	276	9205	0.393	ng/ul	98

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

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