

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071923\
 Data File : BM040961.D
 Acq On : 19 Jul 2023 18:14
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4186

Quant Time: Jul 20 04:59:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 20 04:54:10 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	3147	0.400	ng/ul	0.00
4) Naphthalene-d8	10.680	136	11022	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.527	164	5803	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.282	188	11127	0.400	ng/ul	0.00
17) Chrysene-d12	21.475	240	15532	0.400	ng/ul	0.00
23) Perylene-d12	23.871	264	13285	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	1997	0.404	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.275	152	6118	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.311	212	21308	0.428	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.318	88	2128	0.423	ng/ul#	91
5) Naphthalene	10.730	128	12338	0.406	ng/ul	100
7) 2-Methylnaphthalene	12.352	142	6935	0.388	ng/ul	97
8) 1-Methylnaphthalene	12.566	142	7401	0.415	ng/ul	98
10) Acenaphthylene	14.250	152	9881	0.381	ng/ul	99
11) Acenaphthene	14.587	153	8700	0.397	ng/ul	99
12) Fluorene	15.582	166	9046	0.387	ng/ul	100
14) Pentachlorophenol	16.961	266	995	0.464	ng/ul	99
15) Phenanthrene	17.325	178	13420	0.388	ng/ul	100
16) Anthracene	17.417	178	10661	0.367	ng/ul	99
19) Fluoranthene	19.343	202	26456	0.401	ng/ul	95
20) Pyrene	19.706	202	28891	0.408	ng/ul	96
21) Benzo(a)anthracene	21.457	228	20076	0.382	ng/ul	99
22) Chrysene	21.510	228	21935	0.366	ng/ul	100
24) Benzo(b)fluoranthene	23.141	252	19989	0.389	ng/ul	95
25) Benzo(k)fluoranthene	23.187	252	19874	0.392	ng/ul	95
26) Benzo(a)pyrene	23.769	252	17231	0.372	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.351	276	12150	0.188	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.374	278	9275	0.184	ng/ul	98
29) Benzo(g,h,i)perylene	27.112	276	11134	0.197	ng/ul	99

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

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