

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062821\
 Data File : BM030672.D
 Acq On : 29 Jun 2021 03:19
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4712

Quant Time: Jun 29 04:49:27 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 29 02:23:03 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	2221	0.400	ng/ul	0.00
4) Naphthalene-d8	10.576	136	8610	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.417	164	4662	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.158	188	8611	0.400	ng/ul	0.00
17) Chrysene-d12	21.323	240	7483	0.400	ng/ul	0.00
23) Perylene-d12	23.575	264	7654	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	912	0.377	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.168	152	5105	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.176	212	8578	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	877	0.351	ng/ul#	91
5) Naphthalene	10.625	128	9198	0.375	ng/ul	99
7) 2-Methylnaphthalene	12.244	142	6195	0.375	ng/ul	99
8) 1-Methylnaphthalene	12.460	142	6019	0.375	ng/ul	100
10) Acenaphthylene	14.140	152	7885	0.379	ng/ul	100
11) Acenaphthene	14.481	153	6388	0.379	ng/ul	99
12) Fluorene	15.467	166	6887	0.370	ng/ul	98
14) Pentachlorophenol	16.825	266	862	0.307	ng/ul	96
15) Phenanthrene	17.200	178	10774	0.372	ng/ul	99
16) Anthracene	17.294	178	9339	0.368	ng/ul	99
19) Fluoranthene	19.206	202	12701	0.385	ng/ul	100
20) Pyrene	19.569	202	12681	0.382	ng/ul	99
21) Benzo(a)anthracene	21.306	228	10936	0.367	ng/ul	100
22) Chrysene	21.357	228	12090	0.362	ng/ul	100
24) Benzo(b)fluoranthene	22.897	252	12550	0.370	ng/ul	97
25) Benzo(k)fluoranthene	22.941	252	12620	0.361	ng/ul	98
26) Benzo(a)pyrene	23.479	252	10945	0.380	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.858	276	14559	0.388	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.867	278	11606	0.385	ng/ul	98
29) Benzo(g,h,i)perylene	26.553	276	12592	0.386	ng/ul	99

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

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