

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031668.D
 Acq On : 12 Aug 2021 08:01
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4113

Quant Time: Aug 12 10:40:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 12 10:40:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	1323	0.400	ng/ul	0.00
4) Naphthalene-d8	10.334	136	4677	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.205	164	2389	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	4565	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.157	240	3589	0.400	ng/ul	# 0.00
23) Perylene-d12	23.317	264	3596	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	526	0.358	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.930	152	2837	0.360	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	4597	0.384	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.217	88	539	0.363	ng/ul#	77
5) Naphthalene	10.383	128	5214	0.374	ng/ul	99
7) 2-Methylnaphthalene	12.002	142	3409	0.358	ng/ul	99
8) 1-Methylnaphthalene	12.223	142	3313	0.352	ng/ul	100
10) Acenaphthylene	13.927	152	4143	0.407	ng/ul	100
11) Acenaphthene	14.269	153	3337	0.390	ng/ul	100
12) Fluorene	15.266	166	3618	0.367	ng/ul	98
14) Pentachlorophenol	16.621	266	509	0.348	ng/ul	99
15) Phenanthrene	17.003	178	5426	0.374	ng/ul	99
16) Anthracene	17.096	178	5142	0.379	ng/ul	98
19) Fluoranthene	19.024	202	6025	0.393	ng/ul#	94
20) Pyrene	19.390	202	6280	0.404	ng/ul	96
21) Benzo(a)anthracene	21.142	228	5497	0.379	ng/ul	97
22) Chrysene	21.193	228	5719	0.381	ng/ul	99
24) Benzo(b)fluoranthene	22.672	252	5699	0.352	ng/ul	90
25) Benzo(k)fluoranthene	22.716	252	5824	0.348	ng/ul#	91
26) Benzo(a)pyrene	23.225	252	5149	0.363	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.451	276	6241	0.340	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.468	278	4837	0.327	ng/ul#	91
29) Benzo(g,h,i)perylene	26.103	276	5497	0.351	ng/ul#	92

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

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