

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031268.D
 Acq On : 30 Jul 2021 03:47
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4078

Quant Time: Jul 30 05:12:53 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 : Fri Jul 30 03:45:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	1521	0.400	ng/ul	0.00
4) Naphthalene-d8	10.373	136	5450	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.242	164	3040	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.988	188	5757	0.400	ng/ul	0.00
17) Chrysene-d12	21.178	240	4405	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	4321	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.203	96	605	0.358	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.965	152	3318	0.361	ng/ul	-0.01
18) Fluoranthene-d10	19.020	212	5673	0.386	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	625	0.366	ng/ul#	88
5) Naphthalene	10.423	128	6107	0.376	ng/ul	100
7) 2-Methylnaphthalene	12.042	142	4133	0.373	ng/ul	100
8) 1-Methylnaphthalene	12.267	142	4006	0.366	ng/ul	100
10) Acenaphthylene	13.962	152	4960	0.383	ng/ul	98
11) Acenaphthene	14.303	153	4125	0.378	ng/ul	99
12) Fluorene	15.296	166	4525	0.361	ng/ul	98
14) Pentachlorophenol	16.645	266	679	0.368	ng/ul	99
15) Phenanthrene	17.030	178	6785	0.371	ng/ul	100
16) Anthracene	17.120	178	6385	0.373	ng/ul	99
19) Fluoranthene	19.050	202	7298	0.388	ng/ul	97
20) Pyrene	19.412	202	7492	0.393	ng/ul	97
21) Benzo(a)anthracene	21.161	228	6673	0.375	ng/ul	99
22) Chrysene	21.214	228	6884	0.374	ng/ul	99
24) Benzo(b)fluoranthene	22.698	252	7035	0.361	ng/ul	96
25) Benzo(k)fluoranthene	22.740	252	7031	0.350	ng/ul#	98
26) Benzo(a)pyrene	23.248	252	6269	0.368	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.489	276	8365	0.380	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.504	278	6760	0.381	ng/ul	97
29) Benzo(g,h,i)perylene	26.144	276	7070	0.376	ng/ul	97

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

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