

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031329.D
 Acq On : 01 Aug 2021 02:13
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4083

Quant Time: Aug 02 03:25:50 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 : Mon Aug 02 02:05:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.608	152	1374	0.400	ng/ul	0.00
4) Naphthalene-d8	10.370	136	5230	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.235	164	3022	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.985	188	5673	0.400	ng/ul	0.00
17) Chrysene-d12	21.173	240	4243	0.400	ng/ul	0.00
23) Perylene-d12	23.336	264	4151	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	519	0.340	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.962	152	3222	0.366	ng/ul	0.00
18) Fluoranthene-d10	19.016	212	5398	0.381	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	567	0.368	ng/ul	98
5) Naphthalene	10.419	128	5703	0.366	ng/ul	99
7) 2-Methylnaphthalene	12.038	142	3877	0.364	ng/ul	100
8) 1-Methylnaphthalene	12.259	142	3790	0.360	ng/ul	99
10) Acenaphthylene	13.959	152	4905	0.381	ng/ul	99
11) Acenaphthene	14.300	153	3928	0.363	ng/ul	97
12) Fluorene	15.293	166	4312	0.346	ng/ul	98
14) Pentachlorophenol	16.642	266	624	0.343	ng/ul	98
15) Phenanthrene	17.027	178	6408	0.356	ng/ul	99
16) Anthracene	17.117	178	6039	0.358	ng/ul	99
19) Fluoranthene	19.047	202	6827	0.377	ng/ul	98
20) Pyrene	19.409	202	6846	0.373	ng/ul	98
21) Benzo(a)anthracene	21.159	228	6107	0.356	ng/ul	98
22) Chrysene	21.210	228	6192	0.349	ng/ul	99
24) Benzo(b)fluoranthene	22.691	252	6322	0.338	ng/ul	95
25) Benzo(k)fluoranthene	22.735	252	6368	0.330	ng/ul	97
26) Benzo(a)pyrene	23.241	252	5608	0.342	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.480	276	7430	0.351	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.495	278	6008	0.352	ng/ul	99
29) Benzo(g,h,i)perylene	26.134	276	6346	0.351	ng/ul	96

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

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