

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031680.D
 Acq On : 12 Aug 2021 16:54
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4114

Quant Time: Aug 12 17:27:18 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 11:28:10 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	1230	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	4527	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.208	164	2429	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.964	188	4582	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.158	240	3822	0.400	ng/ul	# 0.00
23) Perylene-d12	23.321	264	3760	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	472	0.346	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	2744	0.360	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	4719	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.220	88	479	0.347	ng/ul#	83
5) Naphthalene	10.382	128	5088	0.377	ng/ul	99
7) 2-Methylnaphthalene	12.006	142	3312	0.360	ng/ul	99
8) 1-Methylnaphthalene	12.226	142	3195	0.351	ng/ul	100
10) Acenaphthylene	13.931	152	4268	0.413	ng/ul	99
11) Acenaphthene	14.273	153	3318	0.381	ng/ul	98
12) Fluorene	15.270	166	3574	0.357	ng/ul	99
14) Pentachlorophenol	16.624	266	534	0.364	ng/ul	99
15) Phenanthrene	17.002	178	5439	0.374	ng/ul	99
16) Anthracene	17.096	178	5224	0.383	ng/ul	98
19) Fluoranthene	19.027	202	6221	0.381	ng/ul	96
20) Pyrene	19.389	202	6392	0.386	ng/ul#	95
21) Benzo(a)anthracene	21.144	228	5725	0.371	ng/ul	97
22) Chrysene	21.195	228	5878	0.368	ng/ul	98
24) Benzo(b)fluoranthene	22.677	252	5866	0.346	ng/ul	88
25) Benzo(k)fluoranthene	22.720	252	5994	0.342	ng/ul#	89
26) Benzo(a)pyrene	23.227	252	5239	0.353	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.460	276	6726	0.351	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.473	278	5328	0.345	ng/ul#	94
29) Benzo(g,h,i)perylene	26.107	276	5655	0.346	ng/ul#	92

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

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