

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100220\
 Data File : BM027577.D
 Acq On : 02 Oct 2020 11:51
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.401

Quant Time: Oct 02 13:07:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 02 13:04:03 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	333	0.40	ng/ul	0.00
2) Naphthalene-d8	10.583	136	875	0.40	ng/ul #	0.00
6) Acenaphthene-d10	14.431	164	671	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.171	188	1770	0.40	ng/ul	0.00
16) Chrysene-d12	21.350	240	2033	0.40	ng/ul	0.00
20) Perylene-d12	23.614	264	2205	0.40	ng/ul #	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.175	152	630	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.198	212	2186	0.45	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.633	128	910	0.36	ng/ul#	87
5) 2-Methylnaphthalene	12.247	142	689	0.37	ng/ul	99
7) Acenaphthylene	14.150	152	1053	0.31	ng/ul#	90
8) Acenaphthene	14.495	153	806	0.33	ng/ul	94
9) Fluorene	15.477	166	1080	0.36	ng/ul#	99
11) Pentachlorophenol	16.820	266	280	0.58	ng/ul	99
12) Phenanthrene	17.213	178	1913	0.35	ng/ul	95
13) Anthracene	17.303	178	1804	0.34	ng/ul#	92
15) Fluoranthene	19.228	202	2725	0.41	ng/ul	99
17) Pyrene	19.590	202	2730	0.29	ng/ul#	96
18) Benzo(a)anthracene	21.335	228	2906	0.36	ng/ul	97
19) Chrysene	21.386	228	2995	0.36	ng/ul	99
21) Benzo(b)fluoranthene	22.934	252	3264	0.37	ng/ul	89
22) Benzo(k)fluoranthene	22.980	252	3256	0.35	ng/ul#	89
23) Benzo(a)pyrene	23.515	252	2933	0.37	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.894	276	4077	0.40	ng/ul#	84
25) Dibenzo(a,h)anthracene	25.911	278	3380	0.40	ng/ul#	90
26) Benzo(g,h,i)perylene	26.594	276	3400	0.39	ng/ul#	95

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

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