

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100720\
 Data File : BM027641.D
 Acq On : 07 Oct 2020 14:04
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.409

Quant Time: Oct 07 14:41:12 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 : Wed Oct 07 11:44:39 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	779	0.40	ng/ul	0.00
2) Naphthalene-d8	10.579	136	3532	0.40	ng/ul #	0.00
6) Acenaphthene-d10	14.423	164	2524	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.168	188	6461	0.40	ng/ul	0.00
16) Chrysene-d12	21.352	240	6833	0.40	ng/ul	0.00
20) Perylene-d12	23.617	264	6785	0.40	ng/ul #	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.167	152	2562	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.194	212	7747	0.44	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.628	128	4045	0.39	ng/ul	99
5) 2-Methylnaphthalene	12.243	142	2940	0.39	ng/ul	100
7) Acenaphthylene	14.143	152	4331	0.34	ng/ul	98
8) Acenaphthene	14.488	153	3410	0.37	ng/ul	96
9) Fluorene	15.473	166	3087	0.27	ng/ul	98
11) Pentachlorophenol	16.817	266	995	0.57	ng/ul	97
12) Phenanthrene	17.209	178	7687	0.39	ng/ul	100
13) Anthracene	17.299	178	7122	0.37	ng/ul	99
15) Fluoranthene	19.224	202	10196	0.43	ng/ul	98
17) Pyrene	19.587	202	10303	0.33	ng/ul#	97
18) Benzo(a)anthracene	21.335	228	10245	0.38	ng/ul	99
19) Chrysene	21.389	228	10916	0.39	ng/ul	100
21) Benzo(b)fluoranthene	22.936	252	11009	0.40	ng/ul	95
22) Benzo(k)fluoranthene	22.980	252	11008	0.39	ng/ul	94
23) Benzo(a)pyrene	23.515	252	9708	0.40	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.897	276	12683	0.40	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.911	278	10332	0.40	ng/ul	97
26) Benzo(g,h,i)perylene	26.592	276	10963	0.41	ng/ul	99

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

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