

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100520\
 Data File : BM027636.D
 Acq On : 06 Oct 2020 13:56
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.407

Quant Time: Oct 06 14:37:15 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 : Mon Oct 05 10:43:07 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	1251	0.40	ng/ul	0.00
2) Naphthalene-d8	10.577	136	3614	0.40	ng/ul #	0.00
6) Acenaphthene-d10	14.425	164	2540	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.166	188	6451	0.40	ng/ul	0.00
16) Chrysene-d12	21.345	240	6850	0.40	ng/ul	0.00
20) Perylene-d12	23.605	264	6992	0.40	ng/ul #	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.164	152	2610	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.192	212	7761	0.44	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.626	128	4112	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.241	142	3021	0.39	ng/ul	98
7) Acenaphthylene	14.140	152	4402	0.34	ng/ul	99
8) Acenaphthene	14.486	153	3423	0.37	ng/ul	98
9) Fluorene	15.471	166	4306	0.38	ng/ul	98
11) Pentachlorophenol	16.815	266	972	0.55	ng/ul	99
12) Phenanthrene	17.208	178	7542	0.38	ng/ul	99
13) Anthracene	17.298	178	7158	0.37	ng/ul	99
15) Fluoranthene	19.222	202	10020	0.42	ng/ul	97
17) Pyrene	19.584	202	10240	0.33	ng/ul#	97
18) Benzo(a)anthracene	21.328	228	10285	0.38	ng/ul	98
19) Chrysene	21.379	228	10778	0.39	ng/ul	99
21) Benzo(b)fluoranthene	22.927	252	11107	0.39	ng/ul	93
22) Benzo(k)fluoranthene	22.970	252	11100	0.38	ng/ul	95
23) Benzo(a)pyrene	23.508	252	10103	0.40	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.883	276	13000	0.40	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.897	278	10622	0.40	ng/ul	95
26) Benzo(g,h,i)perylene	26.578	276	11142	0.41	ng/ul	97

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

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