

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100520\
 Data File : BM027611.D
 Acq On : 05 Oct 2020 18:36
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.405

Quant Time: Oct 05 19:08:51 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 15:37:55 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	1064	0.40	ng/ul	0.00
2) Naphthalene-d8	10.579	136	3036	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.427	164	2056	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.168	188	5262	0.40	ng/ul	0.00
16) Chrysene-d12	21.345	240	5479	0.40	ng/ul	0.00
20) Perylene-d12	23.602	264	5440	0.40	ng/ul	# 0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.167	152	2117	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.194	212	6265	0.43	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.628	128	3482	0.39	ng/ul	97
5) 2-Methylnaphthalene	12.243	142	2491	0.39	ng/ul	98
7) Acenaphthylene	14.143	152	3565	0.34	ng/ul	100
8) Acenaphthene	14.488	153	2827	0.38	ng/ul	96
9) Fluorene	15.473	166	3549	0.38	ng/ul	99
11) Pentachlorophenol	16.817	266	762	0.53	ng/ul	98
12) Phenanthrene	17.209	178	6185	0.38	ng/ul	98
13) Anthracene	17.299	178	5689	0.36	ng/ul	99
15) Fluoranthene	19.220	202	8168	0.42	ng/ul	98
17) Pyrene	19.583	202	8251	0.33	ng/ul	98
18) Benzo(a)anthracene	21.328	228	8167	0.38	ng/ul	99
19) Chrysene	21.379	228	8521	0.38	ng/ul	99
21) Benzo(b)fluoranthene	22.924	252	8738	0.40	ng/ul	92
22) Benzo(k)fluoranthene	22.970	252	8674	0.38	ng/ul	94
23) Benzo(a)pyrene	23.505	252	7676	0.39	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.885	276	10044	0.40	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.897	278	8352	0.40	ng/ul	96
26) Benzo(g,h,i)perylene	26.575	276	8658	0.41	ng/ul	100

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

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