

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100520\
 Data File : BM027627.D
 Acq On : 06 Oct 2020 04:14
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.499

Quant Time: Oct 06 04:48:54 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 : Tue Oct 06 01:23:11 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	1389	0.40	ng/ul	0.00
2) Naphthalene-d8	10.575	136	3940	0.40	ng/ul	# 0.00
6) Acenaphthene-d10	14.423	164	2745	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.164	188	6944	0.40	ng/ul	# 0.00
16) Chrysene-d12	21.340	240	7308	0.40	ng/ul	0.00
20) Perylene-d12	23.600	264	7540	0.40	ng/ul	# 0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.167	152	2816	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.190	212	8359	0.44	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.624	128	4495	0.39	ng/ul	99
5) 2-Methylnaphthalene	12.243	142	3218	0.38	ng/ul	99
7) Acenaphthylene	14.143	152	4771	0.34	ng/ul	99
8) Acenaphthene	14.484	153	3728	0.37	ng/ul	97
9) Fluorene	15.474	166	4718	0.38	ng/ul	99
11) Pentachlorophenol	16.814	266	1109	0.59	ng/ul	99
12) Phenanthrene	17.206	178	8159	0.38	ng/ul	99
13) Anthracene	17.296	178	7715	0.37	ng/ul	99
15) Fluoranthene	19.221	202	10869	0.42	ng/ul	97
17) Pyrene	19.583	202	11025	0.33	ng/ul#	96
18) Benzo(a)anthracene	21.326	228	11100	0.39	ng/ul	99
19) Chrysene	21.377	228	11501	0.39	ng/ul	100
21) Benzo(b)fluoranthene	22.919	252	11941	0.39	ng/ul	94
22) Benzo(k)fluoranthene	22.963	252	11831	0.37	ng/ul	95
23) Benzo(a)pyrene	23.501	252	10802	0.40	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.877	276	14085	0.40	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.892	278	11522	0.40	ng/ul	97
26) Benzo(g,h,i)perylene	26.570	276	11919	0.40	ng/ul	98

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

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