

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111120\
 Data File : BM027791.D
 Acq On : 11 Nov 2020 15:02
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Nov 11 15:35:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 11 11:32:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	926	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	3287	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	1932	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	4047	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	3070	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	2421	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.78	152	2293	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	4463	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	3988	0.380	ng/ul	99
5) 2-Methylnaphthalene	12.85	142	2663	0.364	ng/ul	100
7) Acenaphthylene	14.72	152	3669	0.386	ng/ul	99
8) Acenaphthene	15.05	153	2836	0.389	ng/ul	98
9) Fluorene	16.02	166	3287	0.375	ng/ul	98
11) Pentachlorophenol	17.35	266	546	0.374	ng/ul	99
12) Phenanthrene	17.75	178	5120	0.389	ng/ul	99
13) Anthracene	17.84	178	4892	0.381	ng/ul	100
15) Fluoranthene	19.72	202	5658	0.375	ng/ul	99
17) Pyrene	20.08	202	5646	0.415	ng/ul	99
18) Benzo(a)anthracene	21.79	228	4805	0.392	ng/ul	99
19) Chrysene	21.84	228	4883	0.391	ng/ul	99
21) Benzo(b)fluoranthene	23.50	252	4375	0.395	ng/ul	99
22) Benzo(k)fluoranthene	23.55	252	4310	0.396	ng/ul	99
23) Benzo(a)pyrene	24.14	252	3621	0.372	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.77	276	3953	0.356	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.78	278	3314	0.358	ng/ul	98
26) Benzo(g,h,i)perylene	27.55	276	3363	0.358	ng/ul	98

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

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