

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027987.D
 Acq On : 26 Nov 2020 05:52
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.447

Quant Time: Nov 26 06:25:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 25 10:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	315	0.40	ng/ul	0.00
2) Naphthalene-d8	11.15	136	1195	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.93	164	701	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	1269	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	704	0.40	ng/ul	0.00
20) Perylene-d12	24.17	264	703	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	796	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	1130	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.20	128	1414	0.395	ng/ul#	93
5) 2-Methylnaphthalene	12.79	142	1001	0.411	ng/ul	98
7) Acenaphthylene	14.66	152	1382	0.404	ng/ul	93
8) Acenaphthene	15.00	153	980	0.382	ng/ul	99
9) Fluorene	15.97	166	1125	0.378	ng/ul#	99
11) Pentachlorophenol	17.30	266	155	0.359	ng/ul	97
12) Phenanthrene	17.69	178	1569	0.391	ng/ul	96
13) Anthracene	17.78	178	1518	0.393	ng/ul	94
15) Fluoranthene	19.67	202	1486	0.334	ng/ul	96
17) Pyrene	20.03	202	1436	0.402	ng/ul#	96
18) Benzo(a)anthracene	21.74	228	1128	0.397	ng/ul	93
19) Chrysene	21.79	228	1105	0.399	ng/ul	93
21) Benzo(b)fluoranthene	23.43	252	1085	0.367	ng/ul	93
22) Benzo(k)fluoranthene	23.48	252	1063	0.367	ng/ul	92
23) Benzo(a)pyrene	24.06	252	993	0.384	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	26.66	276	1287	0.460	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.67	278	1069	0.459	ng/ul#	89
26) Benzo(g,h,i)perylene	27.43	276	1090	0.465	ng/ul#	90

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

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