

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027854.D
 Acq On : 20 Nov 2020 15:56
 Operator : JU/CG
 Sample : SSTD3.278
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.278

Quant Time: Nov 20 17:15:19 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 : Fri Nov 20 17:13:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	949	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3743	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2321	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	4909	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	3367	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2499	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	20196	3.20	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	43277	3.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	34215	3.050	ng/ul	97
5) 2-Methylnaphthalene	12.81	142	23796	3.118	ng/ul	99
7) Acenaphthylene	14.68	152	36794	3.248	ng/ul	97
8) Acenaphthene	15.01	153	26511	3.124	ng/ul	99
9) Fluorene	15.98	166	31330	3.182	ng/ul	99
11) Pentachlorophenol	17.31	266	5740	3.433	ng/ul	98
12) Phenanthrene	17.71	178	49017	3.161	ng/ul	97
13) Anthracene	17.80	178	48754	3.259	ng/ul	98
15) Fluoranthene	19.68	202	55364	3.220	ng/ul	99
17) Pyrene	20.04	202	54672	3.197	ng/ul	99
18) Benzo(a)anthracene	21.75	228	42604	3.136	ng/ul	99
19) Chrysene	21.81	228	41624	3.146	ng/ul	98
21) Benzo(b)fluoranthene	23.45	252	33935	3.231	ng/ul	89
22) Benzo(k)fluoranthene	23.50	252	33232	3.231	ng/ul#	90
23) Benzo(a)pyrene	24.08	252	28940	3.152	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.69	276	31433	3.162	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.70	278	26131	3.156	ng/ul#	91
26) Benzo(g,h,i)perylene	27.46	276	26316	3.155	ng/ul	94

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

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