

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027981.D
 Acq On : 26 Nov 2020 01:39
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.446

Quant Time: Nov 26 02:12:49 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 : Thu Nov 26 00:48:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	927	0.40	ng/ul	0.00
2) Naphthalene-d8	11.15	136	3690	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2057	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3774	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	2123	0.40	ng/ul	0.00
20) Perylene-d12	24.17	264	2032	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	2346	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	3389	0.34	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	11.20	128	4450	0.402	ng/ul	99
5) 2-Methylnaphthalene	12.79	142	3026	0.402	ng/ul	99
7) Acenaphthylene	14.66	152	4135	0.412	ng/ul	99
8) Acenaphthene	15.00	153	3021	0.402	ng/ul	99
9) Fluorene	15.97	166	3374	0.387	ng/ul	99
11) Pentachlorophenol	17.30	266	379	0.295	ng/ul	98
12) Phenanthrene	17.69	178	4778	0.401	ng/ul	100
13) Anthracene	17.78	178	4566	0.397	ng/ul	98
15) Fluoranthene	19.67	202	4595	0.348	ng/ul	95
17) Pyrene	20.03	202	4474	0.415	ng/ul#	93
18) Benzo(a)anthracene	21.74	228	3449	0.403	ng/ul	96
19) Chrysene	21.79	228	3361	0.403	ng/ul	98
21) Benzo(b)fluoranthene	23.43	252	3270	0.383	ng/ul	96
22) Benzo(k)fluoranthene	23.48	252	3126	0.374	ng/ul#	94
23) Benzo(a)pyrene	24.06	252	2836	0.380	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	26.66	276	3538	0.438	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.67	278	2911	0.432	ng/ul	96
26) Benzo(g,h,i)perylene	27.43	276	2992	0.441	ng/ul	97

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

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