

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027917.D
 Acq On : 23 Nov 2020 17:24
 Operator : JU/CG
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.440

Quant Time: Nov 23 18:03:58 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 : Mon Nov 23 11:35:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	966	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	3785	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2093	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3731	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2047	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	1912	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2418	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3287	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	4556	0.402	ng/ul	99
5) 2-Methylnaphthalene	12.80	142	3111	0.403	ng/ul	99
7) Acenaphthylene	14.67	152	4150	0.406	ng/ul	99
8) Acenaphthene	15.00	153	3049	0.398	ng/ul	99
9) Fluorene	15.98	166	3372	0.380	ng/ul	100
11) Pentachlorophenol	17.30	266	393	0.309	ng/ul	99
12) Phenanthrene	17.70	178	4701	0.399	ng/ul	100
13) Anthracene	17.79	178	4456	0.392	ng/ul	100
15) Fluoranthene	19.68	202	4424	0.339	ng/ul	97
17) Pyrene	20.04	202	4313	0.415	ng/ul#	95
18) Benzo(a)anthracene	21.75	228	3291	0.398	ng/ul	98
19) Chrysene	21.80	228	3221	0.400	ng/ul	98
21) Benzo(b)fluoranthene	23.45	252	3038	0.378	ng/ul	94
22) Benzo(k)fluoranthene	23.50	252	2925	0.372	ng/ul#	93
23) Benzo(a)pyrene	24.08	252	2656	0.378	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.68	276	3261	0.429	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.69	278	2703	0.427	ng/ul	96
26) Benzo(g,h,i)perylene	27.46	276	2761	0.433	ng/ul	96

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

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