

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027940.D
 Acq On : 24 Nov 2020 13:30
 Operator : JU/CG
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Nov 24 14:15:07 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	8.32	152	1283	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4736	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2260	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3711	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2563	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2557	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	2836	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3609	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	5694	0.401	ng/ul	98
5) 2-Methylnaphthalene	12.80	142	3710	0.384	ng/ul	100
7) Acenaphthylene	14.67	152	4655	0.422	ng/ul	99
8) Acenaphthene	15.00	153	3348	0.405	ng/ul	99
9) Fluorene	15.98	166	3565	0.372	ng/ul	99
11) Pentachlorophenol	17.30	266	449	0.355	ng/ul	97
12) Phenanthrene	17.70	178	4729	0.403	ng/ul	100
13) Anthracene	17.79	178	4653	0.411	ng/ul	98
15) Fluoranthene	19.68	202	4968	0.382	ng/ul	95
17) Pyrene	20.03	202	4896	0.376	ng/ul#	93
18) Benzo(a)anthracene	21.74	228	4179	0.404	ng/ul	95
19) Chrysene	21.80	228	4044	0.401	ng/ul	96
21) Benzo(b)fluoranthene	23.44	252	4102	0.382	ng/ul	95
22) Benzo(k)fluoranthene	23.49	252	3911	0.372	ng/ul#	91
23) Benzo(a)pyrene	24.07	252	3677	0.391	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.67	276	4432	0.436	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.69	278	3605	0.426	ng/ul	97
26) Benzo(g,h,i)perylene	27.44	276	3636	0.426	ng/ul	94

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

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