

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008442.D
 Acq On : 05 Nov 2019 11:55
 Operator : JU
 Sample : SSTDO.102
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.102

Quant Time: Nov 07 10:53:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 10:46:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2247	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	9636	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	5533	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	11562	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	8078	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	7411	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	1560	0.10	ng/ul	0.01
14) Fluoranthene-d10	18.99	212	3561	0.09	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.37	128	2898	0.106	ng/ul	95
5) 2-Methylnaphthalene	12.01	142	1873	0.100	ng/ul	99
7) Acenaphthylene	13.92	152	2670	0.101	ng/ul	96
8) Acenaphthene	14.26	153	2185	0.107	ng/ul	98
9) Fluorene	15.26	166	2349	0.101	ng/ul	96
12) Phenanthrene	16.99	178	3669	0.108	ng/ul	97
13) Anthracene	17.09	178	3066	0.103	ng/ul	96
15) Fluoranthene	19.02	202	4142	0.093	ng/ul	97
17) Pyrene	19.39	202	4369	0.136	ng/ul	98
18) Benzo(a)anthracene	21.15	228	2973	0.101	ng/ul	98
19) Chrysene	21.20	228	3882	0.108	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	2783	0.115	ng/ul	78
22) Benzo(k)fluoranthene	22.73	252	3291	0.120	ng/ul#	82
23) Benzo(a)pyrene	23.24	252	2926	0.115	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.48	276	3338	0.112	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.49	278	2613	0.111	ng/ul#	82
26) Benzo(a,h,i)perylene	26.13	276	3020	0.110	ng/ul	91

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

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