

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003987.D
 Acq On : 16 Dec 2018 08:51
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Dec 17 00:17:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 17 00:13:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3161	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	14736	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8171	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	19597	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19391	0.40	ng/ul	0.01
20) Perylene-d12	23.73	264	16741	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	8598	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	22697	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	15677	0.379	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	11152	0.386	ng/ul	100
7) Acenaphthylene	14.20	152	16758	0.455	ng/ul	99
8) Acenaphthene	14.54	153	12344	0.383	ng/ul	99
9) Fluorene	15.53	166	14050	0.374	ng/ul	99
11) Pentachlorophenol	16.88	266	1923	0.545	ng/ul	96
12) Phenanthrene	17.26	178	23759	0.369	ng/ul	98
13) Anthracene	17.36	178	23918	0.452	ng/ul	99
15) Fluoranthene	19.28	202	29451	0.388	ng/ul	98
17) Pyrene	19.64	202	30107	0.433	ng/ul	98
18) Benzo(a)anthracene	21.39	228	26461	0.432	ng/ul#	88
19) Chrysene	21.44	228	25026	0.357	ng/ul#	89
21) Benzo(b)fluoranthene	23.02	252	24124	0.348	ng/ul#	77
22) Benzo(k)fluoranthene	23.07	252	23912	0.354	ng/ul#	78
23) Benzo(a)pyrene	23.63	252	22509	0.375	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	26.09	276	25930	0.377	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.10	278	21142	0.382	ng/ul#	77
26) Benzo(g,h,i)perylene	26.82	276	21490	0.360	ng/ul#	79

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

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