

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003980.D
 Acq On : 16 Dec 2018 04:44
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Dec 16 23:56:31 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 : Sat Dec 15 23:04:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1622	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8459	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5099	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12816	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	17052	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	15262	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5086	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	15736	0.42	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	8978	0.378	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	6576	0.396	ng/ul	99
7) Acenaphthylene	14.19	152	10063	0.438	ng/ul	99
8) Acenaphthene	14.53	153	7678	0.382	ng/ul	99
9) Fluorene	15.52	166	8885	0.379	ng/ul	100
11) Pentachlorophenol	16.87	266	967	0.419	ng/ul	99
12) Phenanthrene	17.26	178	15492	0.368	ng/ul	99
13) Anthracene	17.35	178	14667	0.424	ng/ul	99
15) Fluoranthene	19.27	202	20835	0.419	ng/ul	96
17) Pyrene	19.64	202	21860	0.357	ng/ul	99
18) Benzo(a)anthracene	21.38	228	22891	0.425	ng/ul	99
19) Chrysene	21.43	228	22348	0.363	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	21961	0.348	ng/ul	96
22) Benzo(k)fluoranthene	23.05	252	21970	0.357	ng/ul#	96
23) Benzo(a)pyrene	23.61	252	20441	0.374	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	26.06	276	22636	0.361	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.08	278	18379	0.364	ng/ul	95
26) Benzo(g,h,i)perylene	26.79	276	18699	0.343	ng/ul	95

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

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