

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002465.D
 Acq On : 23 Aug 2018 14:46
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Aug 23 16:19:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 14:05:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1414	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	5825	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3151	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	7652	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	7816	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	7031	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	3608	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8981	0.40	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.50	128	5984	0.399	ng/ul#	96
5) 2-Methylnaphthalene	12.13	142	4112	0.394	ng/ul	99
7) Acenaphthylene	14.03	152	6015	0.408	ng/ul	96
8) Acenaphthene	14.37	153	4565	0.405	ng/ul	92
9) Fluorene	15.37	166	5234	0.406	ng/ul#	95
11) Pentachlorophenol	16.74	266	287	0.278	ng/ul	90
12) Phenanthrene	17.10	178	9019	0.398	ng/ul#	89
13) Anthracene	17.20	178	7651	0.393	ng/ul#	92
15) Fluoranthene	19.13	202	10556	0.397	ng/ul	97
17) Pyrene	19.49	202	10254	0.385	ng/ul	96
18) Benzo(a)anthracene	21.25	228	9193	0.393	ng/ul#	85
19) Chrysene	21.30	228	10968	0.415	ng/ul	97
21) Benzo(b)fluoranthene	22.82	252	10339	0.409	ng/ul	86
22) Benzo(k)fluoranthene	22.87	252	10517	0.403	ng/ul#	88
23) Benzo(a)pyrene	23.39	252	9277	0.391	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.71	276	9989	0.404	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.72	278	7851	0.403	ng/ul	92
26) Benzo(g,h,i)perylene	26.39	276	8802	0.415	ng/ul	97

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

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