

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092118\
 Data File : BN002785.D
 Acq On : 20 Sep 2018 09:37
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.468

Quant Time: Sep 21 01:15:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 19 17:39:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1291	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	5625	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	3219	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	7474	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	6450	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	5588	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	3526	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	8394	0.44	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.48	128	5879	0.416	ng/ul#	96
5) 2-Methylnaphthalene	12.12	142	4025	0.439	ng/ul	100
7) Acenaphthylene	14.01	152	5461	0.402	ng/ul	98
8) Acenaphthene	14.36	153	4548	0.410	ng/ul	97
9) Fluorene	15.36	166	5195	0.439	ng/ul#	94
11) Pentachlorophenol	16.75	266	202	0.240	ng/ul	92
12) Phenanthrene	17.09	178	8545	0.414	ng/ul#	90
13) Anthracene	17.19	178	7444	0.416	ng/ul#	93
15) Fluoranthene	19.11	202	9730	0.438	ng/ul	96
17) Pyrene	19.48	202	9773	0.465	ng/ul#	97
18) Benzo(a)anthracene	21.24	228	7553	0.424	ng/ul#	85
19) Chrysene	21.29	228	8814	0.424	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	7676	0.416	ng/ul	88
22) Benzo(k)fluoranthene	22.85	252	8251	0.425	ng/ul#	90
23) Benzo(a)pyrene	23.38	252	7221	0.408	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.70	276	7973	0.374	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.70	278	6320	0.370	ng/ul	95
26) Benzo(g,h,i)perylene	26.37	276	6994	0.380	ng/ul	97

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

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