

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003045.D
 Acq On : 10 Oct 2018 01:57
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.498

Quant Time: Oct 10 02:34:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 01:34:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1472	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	6334	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.27	164	3476	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7141	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	4950	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	5104	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.01	152	4131	0.42	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	7639	0.37	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	6894	0.424	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	4714	0.421	ng/ul	99
7) Acenaphthylene	13.99	152	6818	0.429	ng/ul	100
8) Acenaphthene	14.33	153	5255	0.435	ng/ul	100
9) Fluorene	15.34	166	5816	0.421	ng/ul	100
11) Pentachlorophenol	16.74	266	307	0.443	ng/ul	87
12) Phenanthrene	17.07	178	8601	0.432	ng/ul	100
13) Anthracene	17.17	178	7952	0.423	ng/ul	99
15) Fluoranthene	19.09	202	8612	0.369	ng/ul	98
17) Pyrene	19.46	202	8622	0.448	ng/ul	100
18) Benzo(a)anthracene	21.22	228	6956	0.437	ng/ul	99
19) Chrysene	21.27	228	7099	0.439	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	7174	0.390	ng/ul	100
22) Benzo(k)fluoranthene	22.83	252	7552	0.396	ng/ul	99
23) Benzo(a)pyrene	23.36	252	7099	0.431	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.66	276	8876	0.469	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.66	278	7143	0.471	ng/ul	96
26) Benzo(g,h,i)perylene	26.34	276	7487	0.469	ng/ul	97

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

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