

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061318\
 Data File : BN001837.D
 Acq On : 13 Jun 2018 10:54
 Operator : JU/SJ
 Sample : SSTD1.675
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.675

Quant Time: Jun 13 11:32:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:30:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	5841	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	22448	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	12719	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	30027	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	29874	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	27555	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	58235	1.42	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	150460	1.24	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.55	128	93332	1.493	ng/ul	96
5) 2-Methylnaphthalene	12.17	142	67599	1.622	ng/ul	98
7) Acenaphthylene	14.07	152	102224	1.622	ng/ul	95
8) Acenaphthene	14.42	153	74350	1.564	ng/ul	98
9) Fluorene	15.41	166	87577	1.396	ng/ul#	93
11) Pentachlorophenol	16.75	266	9893	2.029	ng/ul	97
12) Phenanthrene	17.14	178	147056	1.656	ng/ul#	89
13) Anthracene	17.23	178	142035	1.661	ng/ul#	92
15) Fluoranthene	19.16	202	176755	1.467	ng/ul	96
17) Pyrene	19.52	202	175375	1.676	ng/ul	97
18) Benzo(a)anthracene	21.28	228	171697	1.748	ng/ul#	85
19) Chrysene	21.33	228	166194	1.694	ng/ul	97
21) Benzo(b)fluoranthene	22.87	252	167311	1.813	ng/ul	83
22) Benzo(k)fluoranthene	22.91	252	167727	1.825	ng/ul#	85
23) Benzo(a)pyrene	23.44	252	156501	1.813	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.78	276	178848	1.859	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.80	278	142993	1.819	ng/ul#	90
26) Benzo(g,h,i)perylene	26.46	276	150178	1.860	ng/ul	96

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

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