

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA N\DATA\BN030218\
 Data File : BN000229.D
 Acq On : 02 Mar 2018 11:02
 Operator : JU/SJ
 Sample : SSTD0.454
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.454

Quant Time: Mar 05 09:37:01 2018
 Quant Method : Z:\HPCHEM1\BNA N\Methods\SOM-EPA-SIM-BN030218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 09:35:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	4954	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	16080	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	7096	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	15960	0.40	ng/ul	0.00
16) Chrysene-d12	21.90	240	10251	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	8241	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	7859	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	13372	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	19121	0.371	ng/ul#	88
5) 2-Methylnaphthalene	12.92	142	12130	0.371	ng/ul	99
7) Acenaphthylene	14.79	152	12934	0.366	ng/ul	96
8) Acenaphthene	15.12	153	12523	0.367	ng/ul	94
9) Fluorene	16.10	166	13558	0.366	ng/ul#	91
11) Pentachlorophenol	17.42	266	998	0.311	ng/ul	97
12) Phenanthrene	17.82	178	23976	0.368	ng/ul#	88
13) Anthracene	17.91	178	17629	0.364	ng/ul#	91
15) Fluoranthene	19.80	202	22836	0.361	ng/ul	62
17) Pyrene	20.16	202	22393	0.371	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	14537	0.361	ng/ul#	84
19) Chrysene	21.94	228	19967	0.365	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	17703	0.357	ng/ul	87
22) Benzo(k)fluoranthene	23.67	252	17002	0.364	ng/ul#	89
23) Benzo(a)pyrene	24.27	252	14090	0.363	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.94	276	15673	0.358	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.95	278	11209	0.355	ng/ul#	80
26) Benzo(g,h,i)perylene	27.72	276	15429	0.358	ng/ul#	82

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

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