

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061318\
 Data File : BN001834.D
 Acq On : 13 Jun 2018 09:08
 Operator : JU/SJ
 Sample : SSTD0.272
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.272

Quant Time: Jun 13 10:33:15 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	6411	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	23679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	13227	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	31824	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	28869	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	25818	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	7530	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	18913	0.15	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.55	128	12307	0.187	ng/ul	96
5) 2-Methylnaphthalene	12.17	142	8709	0.198	ng/ul	98
7) Acenaphthylene	14.07	152	12904	0.197	ng/ul	97
8) Acenaphthene	14.42	153	9599	0.194	ng/ul	97
9) Fluorene	15.41	166	11131	0.171	ng/ul#	93
11) Pentachlorophenol	16.75	266	1044	0.202	ng/ul	98
12) Phenanthrene	17.14	178	19041	0.202	ng/ul#	89
13) Anthracene	17.23	178	17876	0.197	ng/ul#	92
15) Fluoranthene	19.16	202	22004	0.172	ng/ul	96
17) Pyrene	19.53	202	22089	0.218	ng/ul	97
18) Benzo(a)anthracene	21.28	228	19911	0.210	ng/ul#	85
19) Chrysene	21.33	228	19983	0.211	ng/ul	97
21) Benzo(b)fluoranthene	22.87	252	19054	0.220	ng/ul	86
22) Benzo(k)fluoranthene	22.92	252	18676	0.217	ng/ul#	89
23) Benzo(a)pyrene	23.44	252	17172	0.212	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.78	276	17954	0.199	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.80	278	14211	0.193	ng/ul	93
26) Benzo(g,h,i)perylene	26.46	276	15346	0.203	ng/ul	98

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

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