

Data Path : Z:\HPCHEM1\BNA N\DATA\BN031518\
 Data File : BN000435.D
 Acq On : 16 Mar 2018 02:23
 Operator : JU/SJ
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD02008

Quant Time: Mar 16 06:53:55 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN030918MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 16 06:47:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	222385	20.00	ng/ul	0.00
2) Naphthalene-d8	11.30	136	1029526	20.00	ng/ul	0.01
6) Acenaphthene-d10	15.07	164	559352	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.80	188	1161868	20.00	ng/ul	0.01
17) Chrysene-d12	21.92	240	924037	20.00	ng/ul	0.01
22) Perylene-d12	24.42	264	860966	20.00	ng/ul	0.02
System Monitoring Compounds						
7) Acenaphthylene-d8	14.77	160	1131883	20.39	ng/ul	0.00
10) Fluorene-d10	16.05	176	735894	19.57	ng/ul	0.00
14) Anthracene-d10	17.90	188	1077647	19.85	ng/ul	0.00
18) Pyrene-d10	20.15	212	1040838	23.02	ng/ul	0.00
25) Benzo(a)pyrene-d12	24.26	264	791754	20.14	ng/ul	0.02
Target Compounds				Ovalue		
3) Naphthalene	11.35	128	1049780	19.323	ng/ul	100
4) 2-Methylnaphthalene	12.93	142	717424	18.759	ng/ul	97
5) 1-Methylnaphthalene	13.14	142	719285	18.741	ng/ul#	92
8) Acenaphthylene	14.80	152	1110202	19.900	ng/ul	98
9) Acenaphthene	15.14	153	756764	19.490	ng/ul	96
11) Fluorene	16.11	166	829201	19.362	ng/ul	94
13) Phenanthrene	17.84	178	1281919	19.239	ng/ul	99
15) Anthracene	17.93	178	1330575	19.680	ng/ul	99
16) Fluoranthene	19.82	202	1357953	18.862	ng/ul#	93
19) Pyrene	20.18	202	1326809	22.743	ng/ul#	88
20) Benzo(a)anthracene	21.90	228	1154745	20.102	ng/ul	98
21) Chrysene	21.96	228	1066672	19.567	ng/ul	99
23) Benzo(b)fluoranthene	23.65	252	1080409	21.044	ng/ul#	95
24) Benzo(k)fluoranthene	23.71	252	974610	20.280	ng/ul#	96
26) Benzo(a)pyrene	24.31	252	975430	19.963	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	27.01	276	1067175	20.256	ng/ul#	79
28) Dibenzo(a,h)anthracene	27.02	278	899407	20.828	ng/ul#	93
29) Benzo(g,h,i)perylene	27.81	276	890526	20.244	ng/ul#	90

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

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