

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN030518\
 Data File : BN000244.D
 Acq On : 05 Mar 2018 14:34
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.460

Quant Time: Mar 05 15:12:41 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 14:32:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	4710	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	15071	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	6444	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	13928	0.40	ng/ul	0.00
16) Chrysene-d12	21.89	240	8216	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	6538	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	7252	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	11326	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	17929	0.371	ng/ul#	88
5) 2-Methylnaphthalene	12.92	142	11080	0.362	ng/ul	99
7) Acenaphthylene	14.79	152	11462	0.357	ng/ul	97
8) Acenaphthene	15.12	153	11319	0.366	ng/ul	94
9) Fluorene	16.10	166	12126	0.361	ng/ul#	91
11) Pentachlorophenol	17.42	266	820	0.293	ng/ul	98
12) Phenanthrene	17.82	178	20900	0.368	ng/ul#	87
13) Anthracene	17.91	178	14848	0.352	ng/ul#	91
15) Fluoranthene	19.80	202	19336	0.351	ng/ul	62
17) Pyrene	20.15	202	19089	0.395	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	11564	0.359	ng/ul#	83
19) Chrysene	21.93	228	16321	0.372	ng/ul	95
21) Benzo(b)fluoranthene	23.61	252	14684	0.373	ng/ul	87
22) Benzo(k)fluoranthene	23.66	252	13712	0.370	ng/ul#	88
23) Benzo(a)pyrene	24.26	252	11315	0.368	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.91	276	12241	0.352	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.93	278	8585	0.342	ng/ul#	80
26) Benzo(g,h,i)perylene	27.70	276	12221	0.357	ng/ul#	82

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

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