

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053019\
 Data File : BN005955.D
 Acq On : 30 May 2019 12:41
 Operator : JU/SJ
 Sample : SSTD0.289
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.289

Quant Time: May 30 14:05:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 13:56:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	7960	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	29093	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	15762	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	36493	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	23617	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	21828	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	8773	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.18	212	18646	0.16	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	15933	0.207	ng/ul	99
5) 2-Methylnaphthalene	12.19	142	11061	0.210	ng/ul	100
7) Acenaphthylene	14.10	152	17018	0.214	ng/ul	99
8) Acenaphthene	14.45	153	11925	0.212	ng/ul	99
9) Fluorene	15.44	166	13915	0.212	ng/ul	97
11) Pentachlorophenol	16.79	266	1390	0.147	ng/ul	99
12) Phenanthrene	17.18	178	23021	0.232	ng/ul	99
13) Anthracene	17.27	178	21721	0.219	ng/ul	99
15) Fluoranthene	19.21	202	24770	0.180	ng/ul	97
17) Pyrene	19.57	202	25193	0.289	ng/ul	99
18) Benzo(a)anthracene	21.34	228	20571	0.253	ng/ul	98
19) Chrysene	21.39	228	19142	0.191	ng/ul	97
21) Benzo(b)fluoranthene	22.95	252	17371	0.265	ng/ul	84
22) Benzo(k)fluoranthene	22.99	252	17437	0.212	ng/ul#	86
23) Benzo(a)pyrene	23.53	252	16553	0.226	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.93	276	18816	0.229	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.94	278	15139	0.223	ng/ul#	89
26) Benzo(g,h,i)perylene	26.63	276	15983	0.228	ng/ul#	91

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

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