

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110118\
 Data File : BN003373.D
 Acq On : 01 Nov 2018 11:01
 Operator : MJ/SJ
 Sample : SSTD0.427
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.427

Quant Time: Nov 01 11:37:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 11:33:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	11551	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	43949	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	25298	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	61907	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	62110	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	51360	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	28041	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	69019	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.70	128	49019	0.435	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	35267	0.454	ng/ul	100
7) Acenaphthylene	14.21	152	43829	0.379	ng/ul	100
8) Acenaphthene	14.55	153	39130	0.446	ng/ul	100
9) Fluorene	15.54	166	46187	0.459	ng/ul	100
11) Pentachlorophenol	16.87	266	4562	0.759	ng/ul	100
12) Phenanthrene	17.28	178	80295	0.465	ng/ul	100
13) Anthracene	17.36	178	65793	0.404	ng/ul	100
15) Fluoranthene	19.29	202	91005	0.450	ng/ul	100
17) Pyrene	19.65	202	88108	0.365	ng/ul	100
18) Benzo(a)anthracene	21.40	228	77225	0.386	ng/ul	100
19) Chrysene	21.45	228	88061	0.434	ng/ul	100
21) Benzo(b)fluoranthene	23.02	252	86413	0.467	ng/ul	100
22) Benzo(k)fluoranthene	23.07	252	82512	0.430	ng/ul	100
23) Benzo(a)pyrene	23.62	252	72514	0.437	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.08	276	79873	0.419	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.10	278	64983	0.426	ng/ul	100
26) Benzo(g,h,i)perylene	26.81	276	68229	0.425	ng/ul	100

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

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