

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003377.D
 Acq On : 01 Nov 2018 13:30
 Operator : MJ/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV31

Quant Time: Nov 01 14:33:35 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 13:18:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	7795	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	29954	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	16986	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	41633	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	41384	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	32378	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	19091	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	45623	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.70	128	33491	0.399	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	23896	0.395	ng/ul	100
7) Acenaphthylene	14.21	152	28521	0.386	ng/ul	100
8) Acenaphthene	14.55	153	26234	0.398	ng/ul	99
9) Fluorene	15.54	166	30702	0.395	ng/ul	99
11) Pentachlorophenol	16.88	266	2761	0.331	ng/ul	100
12) Phenanthrene	17.28	178	53910	0.395	ng/ul	100
13) Anthracene	17.36	178	42454	0.381	ng/ul	99
15) Fluoranthene	19.29	202	59697	0.389	ng/ul	99
17) Pyrene	19.65	202	58384	0.388	ng/ul	98
18) Benzo(a)anthracene	21.40	228	48921	0.377	ng/ul	99
19) Chrysene	21.45	228	58905	0.399	ng/ul	100
21) Benzo(b)fluoranthene	23.03	252	56980	0.408	ng/ul	99
22) Benzo(k)fluoranthene	23.07	252	52986	0.392	ng/ul	99
23) Benzo(a)pyrene	23.63	252	46216	0.397	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.09	276	49311	0.395	ng/ul	100
25) Dibenzo(a,h)anthracene	26.10	278	40399	0.397	ng/ul	98
26) Benzo(g,h,i)perylene	26.81	276	42426	0.397	ng/ul	99

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

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