

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN002992.D
 Acq On : 08 Oct 2018 18:11
 Operator : MJ/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV93

Quant Time: Oct 08 18:52:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 18:09:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1247	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	5787	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3464	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	8159	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	6811	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4967	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3556	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	9227	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.47	128	5808	0.391	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	4046	0.396	ng/ul	100
7) Acenaphthylene	14.00	152	6176	0.390	ng/ul	100
8) Acenaphthene	14.34	153	4747	0.395	ng/ul	98
9) Fluorene	15.34	166	5474	0.397	ng/ul	99
11) Pentachlorophenol	16.75	266	229	0.289	ng/ul#	89
12) Phenanthrene	17.08	178	8929	0.393	ng/ul	99
13) Anthracene	17.18	178	8352	0.389	ng/ul	99
15) Fluoranthene	19.10	202	10472	0.392	ng/ul	99
17) Pyrene	19.47	202	10651	0.403	ng/ul	99
18) Benzo(a)anthracene	21.23	228	8617	0.393	ng/ul	98
19) Chrysene	21.28	228	8883	0.399	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	6882	0.385	ng/ul	99
22) Benzo(k)fluoranthene	22.84	252	7399	0.399	ng/ul	98
23) Benzo(a)pyrene	23.37	252	6308	0.393	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.67	276	7362	0.400	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.68	278	5879	0.399	ng/ul	100
26) Benzo(g,h,i)perylene	26.35	276	6313	0.406	ng/ul	99

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

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