

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061318\
 Data File : BN001839.D
 Acq On : 13 Jun 2018 12:05
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV77

Quant Time: Jun 13 16:46:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 12:02:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	5768	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	21159	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	11712	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	28074	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	27303	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	25601	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	13189	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	33263	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.55	128	21843	0.401	ng/ul	96
5) 2-Methylnaphthalene	12.17	142	15318	0.393	ng/ul	98
7) Acenaphthylene	14.07	152	22171	0.387	ng/ul	95
8) Acenaphthene	14.42	153	16690	0.396	ng/ul	98
9) Fluorene	15.41	166	19382	0.394	ng/ul#	92
11) Pentachlorophenol	16.74	266	1920	0.361	ng/ul	98
12) Phenanthrene	17.14	178	33372	0.398	ng/ul#	89
13) Anthracene	17.23	178	30932	0.390	ng/ul#	92
15) Fluoranthene	19.16	202	39078	0.393	ng/ul	96
17) Pyrene	19.52	202	38935	0.392	ng/ul	97
18) Benzo(a)anthracene	21.28	228	36701	0.389	ng/ul#	85
19) Chrysene	21.33	228	37632	0.401	ng/ul	96
21) Benzo(b)fluoranthene	22.87	252	36938	0.398	ng/ul	84
22) Benzo(k)fluoranthene	22.92	252	37678	0.407	ng/ul#	87
23) Benzo(a)pyrene	23.44	252	34103	0.395	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.78	276	38667	0.399	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.80	278	31005	0.401	ng/ul#	91
26) Benzo(g,h,i)perylene	26.46	276	33180	0.403	ng/ul	98

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

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