

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062719\
 Data File : BN006602.D
 Acq On : 27 Jun 2019 14:52
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV19

Quant Time: Jun 27 15:31:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 14:44:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1981	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	8171	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	4746	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	10633	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	9747	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	10537	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	5345	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	12888	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	10045	0.428	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	6700	0.423	ng/ul	99
7) Acenaphthylene	14.00	152	10126	0.413	ng/ul	100
8) Acenaphthene	14.34	153	7979	0.423	ng/ul	98
9) Fluorene	15.34	166	8630	0.417	ng/ul	100
11) Pentachlorophenol	16.74	266	518	0.327	ng/ul	98
12) Phenanthrene	17.08	178	14451	0.422	ng/ul	100
13) Anthracene	17.18	178	12045	0.408	ng/ul	100
15) Fluoranthene	19.11	202	17525	0.414	ng/ul	100
17) Pyrene	19.48	202	18458	0.414	ng/ul	100
18) Benzo(a)anthracene	21.27	228	16266	0.406	ng/ul	100
19) Chrysene	21.32	228	18995	0.425	ng/ul	99
21) Benzo(b)fluoranthene	22.87	252	17808	0.412	ng/ul	99
22) Benzo(k)fluoranthene	22.92	252	19401	0.429	ng/ul	100
23) Benzo(a)pyrene	23.45	252	17680	0.418	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.79	276	18424	0.430	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.80	278	14654	0.430	ng/ul	100
26) Benzo(g,h,i)perylene	26.47	276	15626	0.436	ng/ul	99

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

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