

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006986.D
 Acq On : 07 Aug 2019 14:18
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV26

Quant Time: Aug 07 17:40:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	5063	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	19192	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	12108	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	31260	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	29268	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	31748	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	13418	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	39644	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.28	128	21700	0.393	ng/ul	98
5) 2-Methylnaphthalene	11.91	142	16244	0.392	ng/ul	100
7) Acenaphthylene	13.83	152	27124	0.387	ng/ul	99
8) Acenaphthene	14.18	153	18749	0.391	ng/ul	99
9) Fluorene	15.17	166	22935	0.394	ng/ul	100
11) Pentachlorophenol	16.52	266	3035	0.352	ng/ul#	85
12) Phenanthrene	16.90	178	38658	0.388	ng/ul	100
13) Anthracene	16.99	178	39246	0.385	ng/ul	100
15) Fluoranthene	18.93	202	48674	0.393	ng/ul	100
17) Pyrene	19.29	202	49889	0.382	ng/ul	99
18) Benzo(a)anthracene	21.06	228	50267	0.391	ng/ul	99
19) Chrysene	21.11	228	47359	0.393	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	49489	0.384	ng/ul	99
22) Benzo(k)fluoranthene	22.62	252	48014	0.385	ng/ul	99
23) Benzo(a)pyrene	23.11	252	46598	0.385	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.30	276	50987	0.390	ng/ul	100
25) Dibenzo(a,h)anthracene	25.32	278	40887	0.392	ng/ul	100
26) Benzo(g,h,i)perylene	25.93	276	42819	0.391	ng/ul	99

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

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