

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060320\
 Data File : BN010755.D
 Acq On : 03 Jun 2020 16:45
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV10

Quant Time: Jun 03 17:15:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 16:40:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1642	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5336	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2735	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5428	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	4473	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	4138	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	5329	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	9698	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6621	0.420	ng/ul	99
5) 2-Methylnaphthalene	11.99	142	4188	0.424	ng/ul	99
7) Acenaphthylene	13.91	152	5121	0.420	ng/ul	100
8) Acenaphthene	14.25	153	4330	0.424	ng/ul	99
9) Fluorene	15.25	166	4715	0.420	ng/ul	99
11) Pentachlorophenol	16.60	266	573	0.451	ng/ul	96
12) Phenanthrene	16.98	178	7440	0.421	ng/ul	100
13) Anthracene	17.07	178	6039	0.414	ng/ul	99
15) Fluoranthene	19.00	202	8225	0.415	ng/ul	98
17) Pyrene	19.37	202	8130	0.418	ng/ul	99
18) Benzo(a)anthracene	21.13	228	6623	0.417	ng/ul	99
19) Chrysene	21.18	228	8509	0.421	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	7600	0.412	ng/ul	99
22) Benzo(k)fluoranthene	22.71	252	8349	0.428	ng/ul	99
23) Benzo(a)pyrene	23.21	252	6425	0.413	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.42	276	8937	0.419	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.44	278	7340	0.421	ng/ul	99
26) Benzo(g,h,i)perylene	26.06	276	8141	0.413	ng/ul	99

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

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