

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073020\
 Data File : BN011301.D
 Acq On : 30 Jul 2020 16:15
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Jul 30 16:46:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 16:02:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1164	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4015	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2522	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5850	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	5179	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	4711	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	4292	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	11693	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	5034	0.421	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	3434	0.429	ng/ul	97
7) Acenaphthylene	13.82	152	5066	0.421	ng/ul	99
8) Acenaphthene	14.16	153	4054	0.424	ng/ul	99
9) Fluorene	15.16	166	4838	0.430	ng/ul	99
11) Pentachlorophenol	16.52	266	579	0.404	ng/ul	98
12) Phenanthrene	16.90	178	8136	0.420	ng/ul	100
13) Anthracene	16.99	178	7009	0.414	ng/ul	99
15) Fluoranthene	18.93	202	9881	0.413	ng/ul	97
17) Pyrene	19.29	202	9846	0.419	ng/ul	100
18) Benzo(a)anthracene	21.06	228	8228	0.414	ng/ul	100
19) Chrysene	21.11	228	9573	0.420	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	8192	0.417	ng/ul	98
22) Benzo(k)fluoranthene	22.62	252	9607	0.436	ng/ul	99
23) Benzo(a)pyrene	23.11	252	7186	0.420	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.28	276	7484	0.417	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.29	278	5891	0.412	ng/ul	99
26) Benzo(g,h,i)perylene	25.90	276	6660	0.410	ng/ul	99

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

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