

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009337.D
 Acq On : 10 Jan 2020 16:40
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV08

Quant Time: Jan 10 17:10:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 16:45:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1055	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	3620	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2070	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	4513	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3758	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3589	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	2365	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	5446	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	4479	0.411	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	3005	0.402	ng/ul	100
7) Acenaphthylene	13.83	152	4045	0.383	ng/ul	99
8) Acenaphthene	14.18	153	3309	0.390	ng/ul	98
9) Fluorene	15.18	166	3772	0.386	ng/ul	98
11) Pentachlorophenol	16.54	266	428	0.346	ng/ul	98
12) Phenanthrene	16.91	178	6256	0.401	ng/ul	100
13) Anthracene	17.00	178	5296	0.383	ng/ul	99
15) Fluoranthene	18.93	202	7237	0.392	ng/ul	99
17) Pyrene	19.30	202	7324	0.406	ng/ul	98
18) Benzo(a)anthracene	21.06	228	6392	0.388	ng/ul	99
19) Chrysene	21.11	228	7436	0.417	ng/ul	99
21) Benzo(b)fluoranthene	22.57	252	6620	0.416	ng/ul	96
22) Benzo(k)fluoranthene	22.61	252	7659	0.422	ng/ul	97
23) Benzo(a)pyrene	23.11	252	6287	0.394	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.27	276	7097	0.399	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.28	278	5610	0.400	ng/ul	96
26) Benzo(g,h,i)perylene	25.90	276	6347	0.408	ng/ul	100

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

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