

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN020421\
 Data File : BN013571.D
 Acq On : 04 Feb 2021 14:55
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Feb 04 15:31:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 04 14:51:05 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	4575	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	16606	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8691	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18255	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16807	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	12889	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	9933	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	18448	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	18172	0.420	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	12013	0.415	ng/ul	100
7) Acenaphthylene	13.91	152	13225	0.408	ng/ul	99
8) Acenaphthene	14.26	153	11967	0.415	ng/ul	99
9) Fluorene	15.25	166	13474	0.413	ng/ul	100
11) Pentachlorophenol	16.60	266	804	0.362	ng/ul	99
12) Phenanthrene	16.98	178	22263	0.410	ng/ul	100
13) Anthracene	17.07	178	18209	0.404	ng/ul	100
15) Fluoranthene	19.01	202	23441	0.400	ng/ul	100
17) Pyrene	19.37	202	23753	0.393	ng/ul	100
18) Benzo(a)anthracene	21.13	228	22186	0.407	ng/ul	100
19) Chrysene	21.18	228	24656	0.415	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	21602	0.410	ng/ul	98
22) Benzo(k)fluoranthene	22.71	252	22843	0.418	ng/ul	100
23) Benzo(a)pyrene	23.21	252	17418	0.411	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.46	276	23242	0.412	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.47	278	19115	0.413	ng/ul	99
26) Benzo(g,h,i)perylene	26.11	276	20498	0.414	ng/ul	99

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

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