

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022121\
 Data File : BN013700.D
 Acq On : 20 Feb 2021 22:14
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV02

Quant Time: Feb 20 23:19:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN022121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 23:16:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	4014	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	14860	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	8516	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	18500	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17512	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	16281	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	9418	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	21048	0.42	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	17384	0.444	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	11526	0.433	ng/ul	99
7) Acenaphthylene	13.90	152	14235	0.430	ng/ul	100
8) Acenaphthene	14.25	153	12398	0.434	ng/ul	99
9) Fluorene	15.24	166	14299	0.432	ng/ul	99
11) Pentachlorophenol	16.59	266	1672	0.388	ng/ul	98
12) Phenanthrene	16.97	178	24155	0.434	ng/ul	100
13) Anthracene	17.06	178	20877	0.428	ng/ul	100
15) Fluoranthene	19.00	202	27594	0.435	ng/ul	100
17) Pyrene	19.37	202	28182	0.421	ng/ul	97
18) Benzo(a)anthracene	21.12	228	25978	0.431	ng/ul	99
19) Chrysene	21.17	228	27315	0.438	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	27061	0.434	ng/ul	97
22) Benzo(k)fluoranthene	22.70	252	28456	0.444	ng/ul	99
23) Benzo(a)pyrene	23.21	252	25149	0.456	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.45	276	31369	0.441	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.46	278	25657	0.442	ng/ul	98
26) Benzo(g,h,i)perylene	26.10	276	27139	0.442	ng/ul	99

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

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