

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013495.D
 Acq On : 26 Jan 2021 18:50
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Jan 27 03:45:27 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 26 18:47:17 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4845	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	17173	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	8910	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	18605	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	16282	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	16156	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	10856	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	20392	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	20920	0.421	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	13374	0.424	ng/ul	100
7) Acenaphthylene	13.92	152	15124	0.412	ng/ul	100
8) Acenaphthene	14.26	153	13552	0.415	ng/ul	99
9) Fluorene	15.25	166	15106	0.418	ng/ul	100
11) Pentachlorophenol	16.60	266	1092	0.398	ng/ul	98
12) Phenanthrene	16.99	178	25347	0.416	ng/ul	99
13) Anthracene	17.08	178	20691	0.414	ng/ul	100
15) Fluoranthene	19.02	202	27411	0.412	ng/ul	99
17) Pyrene	19.38	202	27306	0.410	ng/ul	98
18) Benzo(a)anthracene	21.14	228	22094	0.411	ng/ul	99
19) Chrysene	21.19	228	27664	0.413	ng/ul	99
21) Benzo(b)fluoranthene	22.68	252	26559	0.392	ng/ul	100
22) Benzo(k)fluoranthene	22.72	252	30234	0.419	ng/ul	100
23) Benzo(a)pyrene	23.23	252	23261	0.404	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.48	276	33021	0.432	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.49	278	26493	0.429	ng/ul	99
26) Benzo(g,h,i)perylene	26.13	276	27347	0.404	ng/ul	100

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

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