

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012848.D
 Acq On : 03 Dec 2020 15:53
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV10

Quant Time: Dec 03 16:24:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 03 15:27:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	665	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.16	136	2376	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	1389	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	3021	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2632	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	2542	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1406	0.40	ng/ul	-0.01
14) Fluoranthene-d10	18.87	212	3284	0.40	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.21	128	2698	0.392	ng/ul	97
5) 2-Methylnaphthalene	11.84	142	1777	0.393	ng/ul	100
7) Acenaphthylene	13.78	152	2582	0.393	ng/ul	98
8) Acenaphthene	14.13	153	2118	0.393	ng/ul	98
9) Fluorene	15.12	166	2448	0.396	ng/ul	97
11) Pentachlorophenol	16.48	266	334	0.382	ng/ul	95
12) Phenanthrene	16.86	178	4073	0.402	ng/ul	99
13) Anthracene	16.96	178	3535	0.410	ng/ul	98
15) Fluoranthene	18.89	202	4718	0.392	ng/ul	99
17) Pyrene	19.26	202	4866	0.391	ng/ul	100
18) Benzo(a)anthracene	21.03	228	4279	0.393	ng/ul	100
19) Chrysene	21.08	228	4698	0.390	ng/ul	99
21) Benzo(b)fluoranthene	22.53	252	4644	0.395	ng/ul	98
22) Benzo(k)fluoranthene	22.57	252	5336	0.411	ng/ul	97
23) Benzo(a)pyrene	23.06	252	4351	0.399	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.22	276	5619	0.392	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.23	278	4488	0.396	ng/ul	95
26) Benzo(g,h,i)perylene	25.84	276	5053	0.390	ng/ul	98

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

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