

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011736.D
 Acq On : 08 Sep 2020 14:41
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV08

Quant Time: Sep 08 15:14:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 14:34:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1693	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	6238	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3399	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7367	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5902	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	5244	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	3456	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	6715	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	6750	0.366	ng/ul	100
5) 2-Methylnaphthalene	12.19	142	4713	0.383	ng/ul	99
7) Acenaphthylene	14.09	152	6157	0.376	ng/ul	99
8) Acenaphthene	14.44	153	4839	0.368	ng/ul	100
9) Fluorene	15.43	166	5568	0.366	ng/ul	95
11) Pentachlorophenol	16.77	266	771	0.347	ng/ul	98
12) Phenanthrene	17.16	178	8983	0.363	ng/ul	99
13) Anthracene	17.25	178	7978	0.362	ng/ul	99
15) Fluoranthene	19.18	202	9915	0.360	ng/ul	99
17) Pyrene	19.55	202	9728	0.371	ng/ul	99
18) Benzo(a)anthracene	21.30	228	8221	0.358	ng/ul	99
19) Chrysene	21.35	228	9608	0.375	ng/ul	99
21) Benzo(b)fluoranthene	22.89	252	9364	0.381	ng/ul	98
22) Benzo(k)fluoranthene	22.93	252	9140	0.363	ng/ul	99
23) Benzo(a)pyrene	23.46	252	7626	0.359	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.83	276	10252	0.368	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	8394	0.369	ng/ul	100
26) Benzo(g,h,i)perylene	26.51	276	8818	0.362	ng/ul	99

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

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