

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060220\
 Data File : BN010746.D
 Acq On : 02 Jun 2020 15:41
 Operator : CG/JU
 Sample : SSTD3.204
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.204

Quant Time: Jun 02 16:17:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 14:30:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2273	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	7662	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	4047	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	8032	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	7644	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	6794	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	39400	3.03	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	78180	3.11	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	74048	3.042	ng/ul	98
5) 2-Methylnaphthalene	11.99	142	48413	2.971	ng/ul	100
7) Acenaphthylene	13.90	152	66059	3.069	ng/ul	98
8) Acenaphthene	14.25	153	50554	3.056	ng/ul	96
9) Fluorene	15.25	166	57590	2.933	ng/ul	100
11) Pentachlorophenol	16.60	266	7686	4.906	ng/ul	99
12) Phenanthrene	16.98	178	89724	3.105	ng/ul	98
13) Anthracene	17.07	178	82237	3.233	ng/ul	98
15) Fluoranthene	19.00	202	105037	2.987	ng/ul	98
17) Pyrene	19.37	202	103131	2.841	ng/ul	99
18) Benzo(a)anthracene	21.13	228	96663	2.936	ng/ul	99
19) Chrysene	21.18	228	102530	2.730	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	97033	3.155	ng/ul	89
22) Benzo(k)fluoranthene	22.71	252	104689	3.226	ng/ul#	89
23) Benzo(a)pyrene	23.21	252	87045	3.265	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.42	276	118407	3.547	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.43	278	97229	3.557	ng/ul	93
26) Benzo(g,h,i)perylene	26.06	276	100962	3.440	ng/ul	95

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

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