

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005490.D
 Acq On : 06 May 2019 18:10
 Operator : JU/SJ
 Sample : SSTD0.259
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.259

Quant Time: May 07 02:50:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 02:47:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1815	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	6648	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.25	164	3779	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9878	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	10951	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	12824	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2023	0.19	ng/ul	0.03
14) Fluoranthene-d10	19.04	212	6208	0.22	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.43	128	3458	0.191	ng/ul#	92
5) 2-Methylnaphthalene	12.08	142	2302	0.180	ng/ul	97
7) Acenaphthylene	13.97	152	3718	0.196	ng/ul	96
8) Acenaphthene	14.31	153	2619	0.187	ng/ul	96
9) Fluorene	15.33	166	3068	0.190	ng/ul#	98
11) Pentachlorophenol	16.68	266	419	0.175	ng/ul	90
12) Phenanthrene	17.05	178	5174	0.173	ng/ul	97
13) Anthracene	17.15	178	5194	0.187	ng/ul	98
15) Fluoranthene	19.07	202	7430	0.211	ng/ul	96
17) Pyrene	19.44	202	8119	0.178	ng/ul	98
18) Benzo(a)anthracene	21.22	228	7271	0.177	ng/ul	98
19) Chrysene	21.27	228	9397	0.216	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	7345	0.156	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	9736	0.200	ng/ul#	94
23) Benzo(a)pyrene	23.35	252	8520	0.196	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.65	276	9451	0.196	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.67	278	7719	0.195	ng/ul	95
26) Benzo(g,h,i)perylene	26.31	276	8059	0.196	ng/ul	95

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

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