

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN053019\
 Data File : BN005956.D
 Acq On : 30 May 2019 13:16
 Operator : JU/SJ
 Sample : SSTD0.490
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.490

Quant Time: May 30 14:03:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 13:56:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	8618	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	31859	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	17065	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	39708	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	26915	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	23877	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	19195	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.18	212	42134	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	34649	0.411	ng/ul	100
5) 2-Methylnaphthalene	12.19	142	24217	0.421	ng/ul	100
7) Acenaphthylene	14.11	152	37557	0.437	ng/ul	100
8) Acenaphthene	14.45	153	26089	0.429	ng/ul	100
9) Fluorene	15.44	166	30441	0.429	ng/ul	100
11) Pentachlorophenol	16.79	266	3382	0.329	ng/ul	100
12) Phenanthrene	17.18	178	50267	0.465	ng/ul	100
13) Anthracene	17.27	178	48068	0.445	ng/ul	100
15) Fluoranthene	19.21	202	55543	0.371	ng/ul	100
17) Pyrene	19.57	202	56250	0.566	ng/ul	100
18) Benzo(a)anthracene	21.34	228	47128	0.508	ng/ul	100
19) Chrysene	21.39	228	44410	0.390	ng/ul	100
21) Benzo(b)fluoranthene	22.95	252	39129	0.545	ng/ul	100
22) Benzo(k)fluoranthene	22.99	252	38843	0.432	ng/ul	100
23) Benzo(a)pyrene	23.53	252	36767	0.458	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.93	276	39783	0.443	ng/ul	100
25) Dibenzo(a,h)anthracene	25.94	278	31654	0.427	ng/ul	100
26) Benzo(g,h,i)perylene	26.63	276	33375	0.436	ng/ul	100

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

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