

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031620\
 Data File : BN010225.D
 Acq On : 16 Mar 2020 13:15
 Operator : CG/JU
 Sample : SSTDO.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.406

Quant Time: Mar 16 14:00:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN031620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 13:58:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1869	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6029	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3534	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	8393	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	7658	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	7770	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	3862	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	9919	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	7206	0.397	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	4825	0.387	ng/ul	100
7) Acenaphthylene	13.97	152	7064	0.392	ng/ul	100
8) Acenaphthene	14.32	153	5454	0.376	ng/ul	100
9) Fluorene	15.31	166	6470	0.388	ng/ul	100
11) Pentachlorophenol	16.65	266	572	0.249	ng/ul	100
12) Phenanthrene	17.04	178	11384	0.393	ng/ul	100
13) Anthracene	17.13	178	9907	0.385	ng/ul	100
15) Fluoranthene	19.06	202	13762	0.401	ng/ul	100
17) Pyrene	19.42	202	13974	0.380	ng/ul	100
18) Benzo(a)anthracene	21.18	228	12354	0.368	ng/ul	100
19) Chrysene	21.23	228	14198	0.391	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	13107	0.380	ng/ul	100
22) Benzo(k)fluoranthene	22.77	252	13768	0.351	ng/ul	100
23) Benzo(a)pyrene	23.27	252	11233	0.326	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.52	276	14090	0.366	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	11644	0.383	ng/ul	100
26) Benzo(g,h,i)perylene	26.17	276	12415	0.368	ng/ul	100

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

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