

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008448.D
 Acq On : 05 Nov 2019 15:48
 Operator : JU
 Sample : SSTDO.404
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.404

Quant Time: Nov 07 10:51:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 10:46:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1337	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5602	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3402	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6506	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	4965	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	4396	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	3496	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	7855	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6317	0.396	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	4194	0.383	ng/ul	100
7) Acenaphthylene	13.91	152	5810	0.359	ng/ul	100
8) Acenaphthene	14.26	153	4759	0.378	ng/ul	100
9) Fluorene	15.25	166	5198	0.363	ng/ul	100
11) Pentachlorophenol	16.62	266	547	0.348	ng/ul	100
12) Phenanthrene	16.99	178	7971	0.418	ng/ul	100
13) Anthracene	17.09	178	6871	0.410	ng/ul	100
15) Fluoranthene	19.02	202	9181	0.365	ng/ul	100
17) Pyrene	19.39	202	9498	0.480	ng/ul	100
18) Benzo(a)anthracene	21.15	228	6748	0.374	ng/ul	100
19) Chrysene	21.20	228	8839	0.399	ng/ul	100
21) Benzo(b)fluoranthene	22.69	252	6699	0.468	ng/ul	100
22) Benzo(k)fluoranthene	22.73	252	7369	0.455	ng/ul	100
23) Benzo(a)pyrene	23.24	252	6620	0.440	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.48	276	7398	0.417	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.50	278	5821	0.415	ng/ul	100
26) Benzo(g,h,i)perylene	26.13	276	6656	0.409	ng/ul	100

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

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