

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101019\
 Data File : BN008060.D
 Acq On : 10 Oct 2019 13:29
 Operator : JU
 Sample : SSTDO.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.403

Quant Time: Oct 10 16:14:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 16:00:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1834	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	8689	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	6041	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	13002	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13512	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	16572	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	6246	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	19272	0.46	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	10554	0.430	ng/ul	100
5) 2-Methylnaphthalene	12.08	142	7221	0.438	ng/ul	100
7) Acenaphthylene	13.99	152	11453	0.341	ng/ul	100
8) Acenaphthene	14.33	153	8965	0.371	ng/ul	100
9) Fluorene	15.32	166	10187	0.390	ng/ul	100
11) Pentachlorophenol	16.68	266	1253	0.428	ng/ul	100
12) Phenanthrene	17.05	178	16334	0.403	ng/ul	100
13) Anthracene	17.15	178	14253	0.373	ng/ul	100
15) Fluoranthene	19.08	202	21538	0.397	ng/ul	100
17) Pyrene	19.44	202	23190	0.351	ng/ul	100
18) Benzo(a)anthracene	21.20	228	20299	0.336	ng/ul	100
19) Chrysene	21.25	228	26428	0.407	ng/ul	100
21) Benzo(b)fluoranthene	22.76	252	21861	0.302	ng/ul	100
22) Benzo(k)fluoranthene	22.80	252	26139	0.345	ng/ul	100
23) Benzo(a)pyrene	23.31	252	23455	0.325	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.59	276	27648	0.342	ng/ul	100
25) Dibenzo(a,h)anthracene	25.60	278	21744	0.340	ng/ul	100
26) Benzo(g,h,i)perylene	26.25	276	25608	0.371	ng/ul	100

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

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