

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009333.D
 Acq On : 10 Jan 2020 14:16
 Operator : JU
 Sample : SSTDO.405
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.405

Quant Time: Jan 10 14:55:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 14:47:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1243	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	4203	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2485	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	5464	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	4576	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	4675	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	2774	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	6716	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	5121	0.402	ng/ul	100
5) 2-Methylnaphthalene	11.91	142	3514	0.365	ng/ul	100
7) Acenaphthylene	13.83	152	4842	0.371	ng/ul	100
8) Acenaphthene	14.18	153	3957	0.386	ng/ul	100
9) Fluorene	15.17	166	4541	0.375	ng/ul	100
11) Pentachlorophenol	16.54	266	573	0.393	ng/ul	100
12) Phenanthrene	16.91	178	7623	0.401	ng/ul	100
13) Anthracene	17.00	178	6675	0.392	ng/ul	100
15) Fluoranthene	18.93	202	8929	0.411	ng/ul	100
17) Pyrene	19.30	202	9131	0.391	ng/ul	100
18) Benzo(a)anthracene	21.05	228	8031	0.395	ng/ul	100
19) Chrysene	21.10	228	9111	0.446	ng/ul	100
21) Benzo(b)fluoranthene	22.57	252	8439	0.387	ng/ul	100
22) Benzo(k)fluoranthene	22.61	252	9609	0.437	ng/ul	100
23) Benzo(a)pyrene	23.10	252	8291	0.433	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.27	276	9284	0.446	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.28	278	7306	0.441	ng/ul	100
26) Benzo(g,h,i)perylene	25.89	276	8213	0.468	ng/ul	100

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

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