

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110118\
 Data File : BN003382.D
 Acq On : 01 Nov 2018 17:14
 Operator : MJ/SJ
 Sample : J5701-21MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40K4MS

Quant Time: Nov 01 18:13:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:18:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10277	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	41116	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	23861	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	59063	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	58564	0.40	ng/ul	0.00
20) Perylene-d12	23.74	264	48331	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.24	152	9961	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	28748	0.17	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) Acenaphthene	14.55	153	13682	0.148	ng/ul	99
12) Phenanthrene	17.27	178	35825	0.185	ng/ul	99
13) Anthracene	17.36	178	3215	0.020	ng/ul#	92
15) Fluoranthene	19.29	202	124870	0.574	ng/ul	99
17) Pyrene	19.65	202	140994	0.662	ng/ul	100
18) Benzo(a)anthracene	21.40	228	44524	0.242	ng/ul	98
19) Chrysene	21.45	228	59727	0.286	ng/ul	98
21) Benzo(b)fluoranthene	23.03	252	105537	0.507	ng/ul	96
22) Benzo(k)fluoranthene	23.03	252	105537	0.523	ng/ul#	96
23) Benzo(a)pyrene	23.63	252	43602	0.251	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.09	276	27490	0.148	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.10	278	6818	0.045	ng/ul#	74
26) Benzo(g,h,i)perylene	26.82	276	20548	0.129	ng/ul	97

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

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