

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101018\
 Data File : BN003056.D
 Acq On : 10 Oct 2018 08:25
 Operator : MJ/SJ
 Sample : J5115-04DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 10 11:27:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1967	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	7640	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3764	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7047	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	5322	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	5640	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.02	152	563	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	891	0.04	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.45	128	642	0.033	ng/ul#	74
5) 2-Methylnaphthalene	12.09	142	478	0.035	ng/ul	92
7) Acenaphthylene	13.99	152	402	0.023	ng/ul#	61
11) Pentachlorophenol	16.73	266	37	0.054	ng/ul#	90
12) Phenanthrene	17.06	178	4993	0.254	ng/ul	98
13) Anthracene	17.15	178	943	0.051	ng/ul#	74
15) Fluoranthene	19.09	202	14589	0.633	ng/ul	100
17) Pyrene	19.45	202	15671	0.758	ng/ul	99
18) Benzo(a)anthracene	21.21	228	10335	0.603	ng/ul	96
19) Chrysene	21.26	228	56230	3.233	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	55826	2.748	ng/ul	96
22) Benzo(k)fluoranthene	22.77	252	55826	2.651	ng/ul	96
23) Benzo(a)pyrene	23.34	252	7307	0.401	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.64	276	14564	0.696	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.64	278	4010	0.239	ng/ul#	76
26) Benzo(g,h,i)perylene	26.32	276	17596	0.998	ng/ul	97

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

