

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082318\
 Data File : BN002466.D
 Acq On : 23 Aug 2018 15:22
 Operator : JU/SJ
 Sample : PB111997BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK97

Quant Time: Aug 23 16:25:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 14:05:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1399	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	5986	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3702	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.15	188	983081	0.40	ng/ul	0.09
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.26
20) Perylene-d12	23.48	264	10392	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.06	152	4082	0.43	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.50	128	210	0.014	ng/ul#	29
5) 2-Methylnaphthalene	12.14	142	111	0.010	ng/ul	87
7) Acenaphthylene	14.02	152	190	0.011	ng/ul#	38
8) Acenaphthene	14.37	153	89	0.007	ng/ul#	83
9) Fluorene	15.37	166	92	0.006	ng/ul#	87
11) Pentachlorophenol	16.84	266	1	0.000	ng/ul#	1
12) Phenanthrene	17.19	178	143	0.000	ng/ul#	41
13) Anthracene	17.40	178	3	0.000	ng/ul#	1
21) Benzo(b)fluoranthene	22.82	252	189	0.005	ng/ul#	1
22) Benzo(k)fluoranthene	22.87	252	174	0.005	ng/ul#	1
23) Benzo(a)pyrene	23.34	252	4574	0.130	ng/ul#	69
24) Indeno(1,2,3-cd)pyrene	25.82	276	124	0.003	ng/ul#	59
25) Dibenzo(a,h)anthracene	25.79	278	138	0.005	ng/ul#	1
26) Benzo(g,h,i)perylene	26.47	276	109	0.003	ng/ul#	1

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

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