

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072219\
 Data File : BN006944.D
 Acq On : 22 Jul 2019 13:54
 Operator : HP/JU
 Sample : K3770-03DL 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2H33DL

Quant Time: Jul 22 14:38:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 23:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1897	0.40	ng/ul	0.02
2) Naphthalene-d8	10.38	136	7410	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.25	164	3947	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8041	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	5661	0.40	ng/ul	0.02
20) Perylene-d12	23.51	264	6309	0.40	ng/ul	0.04

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1792	0.16	ng/ul	0.04
14) Fluoranthene-d10	19.08	212	3860	0.16	ng/ul	0.02

Target Compounds				Ovalue		
3) Naphthalene	10.43	128	10602	0.498	ng/ul	96
7) Acenaphthylene	13.98	152	7308	0.359	ng/ul	98
8) Acenaphthene	14.31	153	6160	0.392	ng/ul	98
9) Fluorene	15.33	166	7141	0.415	ng/ul	98
11) Pentachlorophenol	16.73	266	593	0.495	ng/ul	96
12) Phenanthrene	17.07	178	9391	0.363	ng/ul	98
13) Anthracene	17.18	178	3924	0.176	ng/ul	94
17) Pyrene	19.46	202	23321	0.901	ng/ul	97
18) Benzo(a)anthracene	21.24	228	7149	0.307	ng/ul	100
19) Chrysene	21.29	228	15640	0.603	ng/ul	99
22) Benzo(k)fluoranthene	22.89	252	7125	0.263	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.75	276	10102	0.394	ng/ul#	86
26) Benzo(g,h,i)perylene	26.42	276	12040	0.562	ng/ul	98

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

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