

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070218\
 Data File : BN001930.D
 Acq On : 02 Jul 2018 12:11
 Operator : JU/SJ
 Sample : J3661-03DL 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2D21DL

Quant Time: Jul 03 02:07:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 03 01:51:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	6671	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	26817	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.34	164	16129	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	38143	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	28984	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	23829	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	7847	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	18279	0.16	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	10.53	128	28859	0.418	ng/ul	96
7) Acenaphthylene	14.06	152	25715	0.326	ng/ul	95
8) Acenaphthene	14.40	153	20014	0.345	ng/ul	97
9) Fluorene	15.39	166	28100	0.415	ng/ul#	94
11) Pentachlorophenol	16.74	266	4628	0.641	ng/ul	97
12) Phenanthrene	17.12	178	38001	0.334	ng/ul#	89
13) Anthracene	17.22	178	19650	0.182	ng/ul#	92
17) Pyrene	19.51	202	82080	0.779	ng/ul	97
18) Benzo(a)anthracene	21.27	228	31918	0.318	ng/ul#	85
19) Chrysene	21.32	228	45772	0.459	ng/ul	97
22) Benzo(k)fluoranthene	22.89	252	20228	0.235	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.74	276	28407	0.315	ng/ul#	80
26) Benzo(g,h,i)perylene	26.42	276	30463	0.397	ng/ul	96

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

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