

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070218\
 Data File : BN001928.D
 Acq On : 02 Jul 2018 10:10
 Operator : JU/SJ
 Sample : J3661-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2D21

Quant Time: Jul 03 02:01:23 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	6025	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	24499	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.34	164	14662	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	35217	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	27884	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	23083	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	15099	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	37147	0.35	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.53	128	55443	0.878	ng/ul	96
7) Acenaphthylene	14.06	152	48648	0.679	ng/ul	95
8) Acenaphthene	14.40	153	38772	0.736	ng/ul	98
9) Fluorene	15.39	166	55061	0.894	ng/ul#	93
11) Pentachlorophenol	16.74	266	8807	1.321	ng/ul	96
12) Phenanthrene	17.13	178	75947	0.723	ng/ul#	89
13) Anthracene	17.22	178	38488	0.386	ng/ul#	92
17) Pyrene	19.51	202	171008	1.687	ng/ul	96
18) Benzo(a)anthracene	21.27	228	62541	0.649	ng/ul#	85
19) Chrysene	21.32	228	95081	0.991	ng/ul	97
22) Benzo(k)fluoranthene	22.89	252	42749	0.512	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.74	276	60234	0.690	ng/ul#	80
26) Benzo(g,h,i)perylene	26.41	276	62341	0.840	ng/ul	96

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

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