

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070218\
 Data File : BN001929.D
 Acq On : 02 Jul 2018 11:33
 Operator : JU/SJ
 Sample : J3661-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2D26

Quant Time: Jul 03 01:52:10 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	7.71	152	6163	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	26491	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.34	164	15991	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	38857	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	31295	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	25216	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	12805	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	40899	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	16389	0.240	ng/ul	97
5) 2-Methylnaphthalene	12.15	142	11367	0.233	ng/ul	98
7) Acenaphthylene	14.06	152	35970	0.460	ng/ul	95
9) Fluorene	15.39	166	21072	0.314	ng/ul#	94
11) Pentachlorophenol	16.74	266	1728	0.235	ng/ul	97
12) Phenanthrene	17.12	178	70313	0.607	ng/ul#	89
13) Anthracene	17.22	178	21729	0.198	ng/ul#	92
15) Fluoranthene	19.15	202	53103	0.386	ng/ul	96
17) Pyrene	19.51	202	61088	0.537	ng/ul	97
18) Benzo(a)anthracene	21.27	228	68945	0.637	ng/ul#	85
19) Chrysene	21.32	228	38094	0.354	ng/ul	97
21) Benzo(b)fluoranthene	22.85	252	26206	0.286	ng/ul	90
22) Benzo(k)fluoranthene	22.89	252	45653	0.500	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.74	276	32996	0.346	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.76	278	28916	0.380	ng/ul#	91
26) Benzo(a,h,i)perylene	26.42	276	42107	0.519	ng/ul	96

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

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