

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021861.D
 Acq On : 06 Aug 2019 10:25
 Operator : HP/JU
 Sample : K3890-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A52U7

Quant Time: Aug 06 11:22:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 06 01:29:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	598	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	3113	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1846	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	4073	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3892	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4442	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	1625	0.32	ng/ul	0.03
14) Fluoranthene-d10	19.08	212	4509	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	6354	0.723	ng/ul	97
7) Acenaphthylene	14.01	152	6302	0.732	ng/ul	96
8) Acenaphthene	14.36	153	4794	0.681	ng/ul	98
9) Fluorene	15.36	166	6119	0.782	ng/ul	98
11) Pentachlorophenol	16.71	266	835	1.111	ng/ul	90
12) Phenanthrene	17.09	178	7586	0.640	ng/ul	98
13) Anthracene	17.19	178	4302	0.426	ng/ul	98
17) Pyrene	19.47	202	21763	1.361	ng/ul	99
18) Benzo(a)anthracene	21.23	228	9677	0.695	ng/ul	99
19) Chrysene	21.29	228	15590	0.871	ng/ul	99
22) Benzo(k)fluoranthene	22.85	252	7126	0.420	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.69	276	11422	0.628	ng/ul#	92
26) Benzo(g,h,i)perylene	26.36	276	12785	0.781	ng/ul	98

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

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