

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081519\
 Data File : BM022150.D
 Acq On : 15 Aug 2019 19:10
 Operator : HP/JU
 Sample : K4242-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG6

Quant Time: Aug 16 14:12:26 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 : Wed Aug 14 13:53:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	663	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.37	136	2770	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.25	164	1563	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	213569	0.40	ng/ul	0.07
16) Chrysene-d12	21.36	240	160	0.40	ng/ul	0.13
20) Perylene-d12	0.00	264	0	0.00	ng/ul	-23.44
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	1274	0.28	ng/ul	-0.03
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds						
					Qvalue	
3) Naphthalene	10.42	128	3086	0.394	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	2158	0.414	ng/ul	98
7) Acenaphthylene	13.97	152	1034	0.142	ng/ul#	79
8) Acenaphthene	14.31	153	21020	3.524	ng/ul	96
9) Fluorene	15.31	166	32601	4.921	ng/ul#	97
12) Phenanthrene	17.14	178	118724	0.191	ng/ul	96
17) Pyrene	19.44	202	458293	697.036	ng/ul	96
18) Benzo(a)anthracene	21.25	228	218101	381.089	ng/ul	97
19) Chrysene	21.25	228	218101	296.319	ng/ul	99

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

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