

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081519\
 Data File : BM022154.D
 Acq On : 15 Aug 2019 21:36
 Operator : HP/JU
 Sample : K4242-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AH1

Quant Time: Aug 16 14:12:43 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	621	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.37	136	2761	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.25	164	1595	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	195807	0.40	ng/ul	0.07
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.23
20) Perylene-d12	23.43	264	4323	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	1068	0.23	ng/ul	-0.02
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds						
					Qvalue	
7) Acenaphthylene	13.97	152	150	0.020	ng/ul#	44
21) Benzo(b)fluoranthene	22.77	252	1109	0.076	ng/ul#	20
22) Benzo(k)fluoranthene	22.77	252	1091	0.066	ng/ul#	20
23) Benzo(a)pyrene	23.32	252	407	0.027	ng/ul#	3
24) Indeno(1,2,3-cd)pyrene	25.65	276	409	0.023	ng/ul#	98
26) Benzo(g,h,i)perylene	26.33	276	391	0.025	ng/ul#	22

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

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