

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081519\
 Data File : BM022149.D
 Acq On : 15 Aug 2019 18:33
 Operator : HP/JU
 Sample : K4242-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AF5

Quant Time: Aug 16 14:12:22 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	699	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.37	136	3028	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.25	164	1789	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	282536	0.40	ng/ul	0.07
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.23
20) Perylene-d12	23.42	264	4557	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	1562	0.31	ng/ul	-0.03
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
3) Naphthalene	10.42	128	440	0.051	ng/ul#	68
5) 2-Methylnaphthalene	12.06	142	409	0.072	ng/ul	96
7) Acenaphthylene	13.97	152	4689	0.562	ng/ul#	82
8) Acenaphthene	14.31	153	575	0.084	ng/ul	95
9) Fluorene	15.31	166	1577	0.208	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	69352	4.513	ng/ul	85
22) Benzo(k)fluoranthene	22.76	252	69352	3.988	ng/ul#	85
23) Benzo(a)pyrene	23.32	252	35590	2.277	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.62	276	24101	1.292	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.62	278	6412	0.425	ng/ul#	88
26) Benzo(g,h,i)perylene	26.30	276	18713	1.114	ng/ul	99

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

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