

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081519\
 Data File : BM022152.D
 Acq On : 15 Aug 2019 20:23
 Operator : HP/JU
 Sample : K4242-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ1

Quant Time: Aug 16 14:12:34 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	7.59	152	669	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.37	136	2823	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.25	164	1628	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	300449	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	4571	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	1614	0.35	ng/ul	-0.02
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
7) Acenaphthylene	13.97	152	665	0.088	ng/ul#	74
8) Acenaphthene	14.31	153	135	0.022	ng/ul#	86
9) Fluorene	15.31	166	219	0.032	ng/ul#	90
21) Benzo(b)fluoranthene	22.76	252	8056	0.523	ng/ul	91
22) Benzo(k)fluoranthene	22.76	252	8056	0.462	ng/ul#	91
23) Benzo(a)pyrene	23.32	252	4181	0.267	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.64	276	3123	0.167	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.64	278	759	0.050	ng/ul#	12
26) Benzo(g,h,i)perylene	26.32	276	2805	0.166	ng/ul#	80

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

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