

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032416\
 Data File : BM004754.D
 Acq On : 24 Mar 2016 13:27
 Operator : UM/SJ
 Sample : H1909-22
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4T56

Quant Time: Apr 05 18:23:55 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 16:23:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1180	0.40	ng/ul	0.00
4) Naphthalene-d8	10.62	136	4678	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	2317	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	5074	0.40	ng/ul	0.00
18) Chrysene-d12	21.40	240	4544	0.40	ng/ul	-0.01
22) Perylene-d12	23.69	264	3828	0.40	ng/ul	-0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.22	152	1859	0.35	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	4789	0.47	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.67	128	5921	0.54	ng/ul#	96
7) 2-Methylnaphthalene	12.28	142	4070	0.58	ng/ul	98
9) Acenaphthylene	14.18	152	5039	0.59	ng/ul	96
10) Acenaphthene	14.52	153	5363	0.70	ng/ul	87
13) Pentachlorophenol	16.87	266	528	0.78	ng/ul	95
15) Anthracene	17.33	178	6358	0.56	ng/ul#	94
19) Pyrene	19.63	202	14620	0.92	ng/ul	98
23) Benzo(b)fluoranthene	22.99	252	10975	0.80	ng/ul	94
24) Benzo(k)fluoranthene	23.03	252	16858	1.22	ng/ul#	92

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

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