

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041116\
 Data File : BM004935.D
 Acq On : 11 Apr 2016 13:07
 Operator : UM/SJ
 Sample : H2275-11RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9T84RE

Quant Time: Apr 12 01:30:04 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 12 01:10:05 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1263	0.40	ng/ul	0.00
4) Naphthalene-d8	10.61	136	5466	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.45	164	3160	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.19	188	6857	0.40	ng/ul	0.00
18) Chrysene-d12	21.38	240	7055	0.40	ng/ul	-0.01
22) Perylene-d12	23.67	264	6513	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	632	1.11	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	874	0.11	ng/ul	0.00
16) Fluoranthene-d10	19.23	212	2437	0.15	ng/ul	0.00
Target Compounds				Ovalue		
11) Fluorene	15.50	166	227	0.02	ng/ul#	66
14) Phenanthrene	17.23	178	773	0.04	ng/ul#	75
17) Fluoranthene	19.26	202	626	0.03	ng/ul	71
20) Benzo(a)anthracene	21.37	228	391	0.02	ng/ul#	67
21) Chrysene	21.42	228	390	0.02	ng/ul#	68

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

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