

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041116\
 Data File : BM004934.D
 Acq On : 11 Apr 2016 12:31
 Operator : UM/SJ
 Sample : H2275-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9T75

Manual Integrations
 APPROVED

Sohil
 4/12/2016 5:27:04 PM

Quant Time: Apr 12 01:25:31 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	1303	0.40	ng/ul	0.00
4) Naphthalene-d8	10.61	136	5698	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.45	164	3345	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.19	188	6963m	0.40	ng/ul	0.00
18) Chrysene-d12	21.39	240	7131	0.40	ng/ul	0.00
22) Perylene-d12	23.67	264	5255m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	1648	2.81	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	2182	0.27	ng/ul	0.00
16) Fluoranthene-d10	19.23	212	5827	0.35	ng/ul	0.00
Target Compounds				Ovalue		
7) 2-Methylnaphthalene	12.28	142	268	0.03	ng/ul	99
9) Acenaphthylene	14.16	152	451	0.03	ng/ul#	68
10) Acenaphthene	14.50	153	266	0.02	ng/ul#	76
11) Fluorene	15.50	166	1025m	0.08	ng/ul	
13) Pentachlorophenol	16.85	266	186	0.11	ng/ul	91
14) Phenanthrene	17.23	178	2736m	0.14	ng/ul	
15) Anthracene	17.33	178	2425	0.13	ng/ul#	91
17) Fluoranthene	19.25	202	3519	0.16	ng/ul	93
19) Pyrene	19.62	202	4457	0.18	ng/ul#	95
20) Benzo(a)anthracene	21.37	228	2040	0.09	ng/ul#	93
21) Chrysene	21.42	228	1815	0.09	ng/ul	94
23) Benzo(b)fluoranthene	22.98	252	2025m	0.11	ng/ul	
24) Benzo(k)fluoranthene	23.02	252	1444m	0.09	ng/ul	
26) Indeno(1,2,3-cd)pyrene	26.00	276	1222m	0.08	ng/ul	
27) Dibenzo(a,h)anthracene	26.01	278	883m	0.07	ng/ul	
28) Benzo(a,h,i)perylene	26.70	276	923m	0.07	ng/ul	

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

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