

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120915\
 Data File : BM003454.D
 Acq On : 10 Dec 2015 16:08
 Operator : UM/IZ
 Sample : G4713-02
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TEC-17

Manual Integrations
 APPROVED

MMdadoda
 12/10/2015 5:57:58 PM

Quant Time: Dec 10 17:55:46 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM120915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 18:52:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	82641	20.00	ng	0.00
21) Naphthalene-d8	10.52	136	358871	20.00	ng	0.00
38) Acenaphthene-d10	14.37	164	185270	20.00	ng	0.00
63) Phenanthrene-d10	17.12	188	347072	20.00	ng	0.00
75) Chrysene-d12	21.32	240	297628	20.00	ng	0.00
86) Perylene-d12	23.58	264	301440	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	220311	47.40	ng	0.00
7) Phenol-d6	6.89	99	186719	28.93	ng	0.00
23) Nitrobenzene-d5	8.88	82	566571	97.89	ng	0.00
41) 2,4,6-Tribromophenol	15.87	330	181040	97.84	ng	0.00
44) 2-Fluorobiphenyl	12.99	172	1304834	95.95	ng	0.00
78) Terphenyl-d14	19.76	244	1100993	87.24	ng	0.00
Target Compounds						
9) Benzaldehyde	6.86	77	14114	4.46	ng	# 87
10) Phenol	6.92	94	103770	15.15	ng	# 72
20) 3+4-Methylphenols	8.49	107	16280	2.46	ng	98
32) Benzoic acid	9.85	122	140179m	42.33	ng	

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

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