

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
 Data File : BF081023.D
 Acq On : 18 Aug 2015 1:39
 Operator : UM/IZ
 Sample : G3252-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-01-017-A

Manual Integrations
 APPROVED

MMDadoda
 8/18/2015 5:12:52 PM

Quant Time: Aug 18 06:46:03 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 17:57:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	72446	20.00	ng	0.00
21) Naphthalene-d8	7.91	136	257908	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	122518	20.00	ng	0.00
63) Phenanthrene-d10	11.13	188	247768	20.00	ng	0.00
75) Chrysene-d12	13.75	240	161203	20.00	ng	0.00
86) Perylene-d12	15.10	264	143944	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	572581	118.88	ng	0.02
7) Phenol-d6	6.27	99	740825	117.88	ng	0.01
23) Nitrobenzene-d5	7.20	82	556833	98.63	ng	0.01
41) 2,4,6-Tribromophenol	10.44	330	128769	121.25	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	660796	70.67	ng	0.00
78) Terphenyl-d14	12.72	244	565415	75.19	ng	0.01
Target Compounds						
49) Dimethylphthalate	9.39	163	170577	16.91	ng	Qvalue 98
74) Fluoranthene	12.33	202	60706	3.83	ng	95
77) Pyrene	12.56	202	52081	3.97	ng	97
80) Benzo(a)anthracene	13.74	228	29511	2.73	ng	94
82) Chrysene	13.77	228	21623m	2.22	ng	
85) Indeno(1,2,3-cd)pyrene	16.33	276	14531	2.12	ng	# 81
88) Benzo(k)fluoranthene	14.73	252	27440m	2.89	ng	
89) Benzo(a)pyrene	15.04	252	22900	2.58	ng	# 88

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

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