

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\
 Data File : BG021207.D
 Acq On : 2 Mar 2016 6:20
 Operator : UM/SJ
 Sample : H1557-01RE 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P-09RE

Manual Integrations
 APPROVED

Sohil
 3/2/2016 9:52:28 AM

Quant Time: Mar 02 12:48:51 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 29 17:40:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	13943	20.00	ng	0.00
21) Naphthalene-d8	10.96	136	66559	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	36227	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	74943	20.00	ng	0.00
75) Chrysene-d12	21.76	240	81478	20.00	ng	0.00
86) Perylene-d12	25.04	264	61186	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	5883	6.71	ng	0.00
7) Phenol-d6	7.29	99	9384	6.51	ng	0.00
23) Nitrobenzene-d5	9.31	82	7237	4.60	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	2372	6.29	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	11896	4.66	ng	0.00
78) Terphenyl-d14	20.08	244	15159	4.51	ng	-0.01
Target Compounds						
74) Fluoranthene	19.53	202	30190	6.35	ng	99
77) Pyrene	19.89	202	30379	6.25	ng	95
80) Benzo(a)anthracene	21.74	228	22338	4.63	ng	# 95
82) Chrysene	21.80	228	22455m	4.91	ng	
87) Benzo(b)fluoranthene	24.00	252	34766	9.03	ng	# 89
88) Benzo(k)fluoranthene	24.05	252	14376m	3.73	ng	

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

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