

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051216\
 Data File : BF086954.D
 Acq On : 13 May 2016 2:44
 Operator : UM/SJ
 Sample : H2968-07 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUPB

Manual Integrations
 APPROVED

sohil
 5/13/2016 8:08:20 PM

Quant Time: May 13 04:26:52 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 13 00:58:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	159654	20.00	ng	-0.01
21) Naphthalene-d8	7.91	136	622811	20.00	ng	-0.01
38) Acenaphthene-d10	9.66	164	268151	20.00	ng	-0.01
63) Phenanthrene-d10	11.13	188	408938	20.00	ng	-0.01
75) Chrysene-d12	13.76	240	391339	20.00	ng	-0.01
86) Perylene-d12	15.13	264	303453	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	153859	16.32	ng	0.00
7) Phenol-d6	6.28	99	219952	17.46	ng	-0.01
23) Nitrobenzene-d5	7.19	82	125793	10.14	ng	-0.02
41) 2,4,6-Tribromophenol	10.44	330	31546	11.11	ng	-0.01
44) 2-Fluorobiphenyl	8.98	172	199635	12.51	ng	-0.01
78) Terphenyl-d14	12.72	244	142636	7.73	ng	0.00
Target Compounds						
74) Fluoranthene	12.34	202	162047	7.22	ng	Qvalue 93
77) Pyrene	12.56	202	132874	4.43	ng	96
80) Benzo(a)anthracene	13.75	228	92830	4.23	ng	93
82) Chrysene	13.78	228	61863m	2.77	ng	
87) Benzo(b)fluoranthene	14.75	252	77503m	3.93	ng	
88) Benzo(k)fluoranthene	14.77	252	33073m	2.05	ng	
89) Benzo(a)pyrene	15.07	252	50502	2.97	ng	# 92

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

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