

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
 Data File : BF085687.D
 Acq On : 23 Mar 2016 5:32
 Operator : UM/SJ
 Sample : H1858-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-12A

Manual Integrations
 APPROVED

Sohil
 3/23/2016 1:30:22 PM

Quant Time: Mar 23 06:52:05 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	137701	20.00	ng	-0.02
21) Naphthalene-d8	8.12	136	569625	20.00	ng	-0.02
38) Acenaphthene-d10	9.86	164	240007	20.00	ng	-0.03
63) Phenanthrene-d10	11.35	188	432134	20.00	ng	-0.03
75) Chrysene-d12	13.98	240	251011	20.00	ng	-0.03
86) Perylene-d12	15.40	264	241305	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1213167	147.98	ng	-0.01
7) Phenol-d6	6.47	99	1567751	149.10	ng	-0.02
23) Nitrobenzene-d5	7.40	82	973432	99.19	ng	-0.03
41) 2,4,6-Tribromophenol	10.66	330	419602	178.48	ng	-0.02
44) 2-Fluorobiphenyl	9.20	172	1702763	147.43	ng	-0.02
78) Terphenyl-d14	12.94	244	1250022	119.84	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.59	163	364043	20.10	ng	99
74) Fluoranthene	12.56	202	61942	2.60	ng	90
77) Pyrene	12.79	202	63972	3.55	ng	99
80) Benzo(a)anthracene	13.97	228	39338	2.70	ng	94
82) Chrysene	14.00	228	29168m	2.08	ng	
87) Benzo(b)fluoranthene	15.00	252	55622m	3.53	ng	

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

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