

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
 Data File : BF088546.D
 Acq On : 30 Jun 2016 23:23
 Operator : UM/SJ
 Sample : H3629-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOCATION-5D(9-11)

Manual Integrations

APPROVED
 SOHIL
 7/1/2016 8:05:27 AM

Quant Time: Jul 01 02:19:33 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 30 18:01:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	175783	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	748576	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	352774	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	631620	20.00	ng	0.00
75) Chrysene-d12	13.95	240	482892	20.00	ng	0.00
86) Perylene-d12	15.35	264	387526	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1236925	105.46	ng	0.01
7) Phenol-d6	6.46	99	1934109	127.62	ng	0.01
23) Nitrobenzene-d5	7.38	82	1195128	83.67	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	410834	125.41	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	1454008	65.10	ng	0.00
78) Terphenyl-d14	12.91	244	1208523	55.74	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	275429	10.96	ng	99
70) Phenanthrene	11.35	178	163395	4.73	ng	98
74) Fluoranthene	12.53	202	203187	5.80	ng	95
77) Pyrene	12.75	202	208995	5.10	ng	99
80) Benzo(a)anthracene	13.94	228	78677	2.53	ng	99
82) Chrysene	13.98	228	71623m	2.33	ng	
83) Bis(2-ethylhexyl)phthalate	13.94	149	52734	2.53	ng	# 98
87) Benzo(b)fluoranthene	14.95	252	61519m	2.53	ng	
89) Benzo(a)pyrene	15.29	252	51067m	2.48	ng	

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

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