

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081916\
 Data File : BF089803.D
 Acq On : 19 Aug 2016 18:40
 Operator : UM/SJ
 Sample : H4463-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STONE-1

Manual Integrations
 APPROVED

sohil
 8/22/2016 3:48:07 PM

Quant Time: Aug 22 01:47:59 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 01:47:27 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	567713	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	2064596	20.00	ng	-0.01
38) Acenaphthene-d10	9.71	164	973239	20.00	ng	-0.01
63) Phenanthrene-d10	11.19	188	1665029	20.00	ng	-0.01
75) Chrysene-d12	13.83	240	1118739	20.00	ng	-0.07
86) Perylene-d12	15.26	264	984867	20.00	ng	-0.19
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	3421064	103.34	ng	0.01
7) Phenol-d6	6.32	99	3556098	87.51	ng	0.00
23) Nitrobenzene-d5	7.24	82	2171465	65.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.50	330	897538	97.04	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	4380151	79.10	ng	0.00
78) Terphenyl-d14	12.78	244	3230650	65.33	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.44	163	1215697	17.74	ng	99
70) Phenanthrene	11.21	178	440914	5.24	ng	98
74) Fluoranthene	12.39	202	447528	5.59	ng	93
77) Pyrene	12.62	202	441376	4.84	ng	97
80) Benzo(a)anthracene	13.82	228	165071m	2.58	ng	
82) Chrysene	13.85	228	148838m	2.71	ng	
87) Benzo(b)fluoranthene	14.88	252	136197m	2.51	ng	
89) Benzo(a)pyrene	15.20	252	113636	2.22	ng	# 96

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

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