

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032917\
 Data File : BF094051.D
 Acq On : 30 Mar 2017 8:07
 Operator : SJ/MA
 Sample : I2431-04
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-T12-SSW

Manual Integrations
 APPROVED

mohammad
 3/30/2017 2:32:27 PM

Quant Time: Mar 30 08:29:37 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 27 16:10:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.92	152	168821	20.00	ng	-0.03
21) Naphthalene-d8	7.43	136	654710	20.00	ng	-0.03
38) Acenaphthene-d10	9.57	164	281232	20.00	ng	-0.03
63) Phenanthrene-d10	11.39	188	441189	20.00	ng	-0.03
75) Chrysene-d12	14.64	240	313350	20.00	ng	-0.04
86) Perylene-d12	16.28	264	253288	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	4.46	112	1303911	130.45	ng	-0.01
7) Phenol-d6	5.53	99	1622853	131.24	ng	-0.02
23) Nitrobenzene-d5	6.57	82	858318	74.82	ng	-0.03
41) 2,4,6-Tribromophenol	10.55	330	262487	118.11	ng	-0.03
44) 2-Fluorobiphenyl	8.75	172	1674544	90.82	ng	-0.03
78) Terphenyl-d14	13.37	244	1054548	75.44	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.25	163	108084	5.41	ng	98
70) Phenanthrene	11.42	178	56354	2.30	ng	98
74) Fluoranthene	12.89	202	104283	4.15	ng	96
77) Pyrene	13.16	202	98530	4.33	ng	97
80) Benzo(a)anthracene	14.63	228	51890	2.84	ng	94
82) Chrysene	14.67	228	45562	2.69	ng	94
87) Benzo(b)fluoranthene	15.87	252	68463m	4.41	ng	
89) Benzo(a)pyrene	16.22	252	41718	2.92	ng	# 95

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

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