

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071317\
 Data File : BF096745.D
 Acq On : 13 Jul 2017 23:30
 Operator : SJ/JU
 Sample : I4146-03 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-001-B

Manual Integrations
 APPROVED

mohammad
 7/17/2017 8:11:52 AM

Quant Time: Jul 14 12:41:20 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 14:49:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	94666	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	346380	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	124451	20.00	ng	0.00
63) Phenanthrene-d10	11.15	188	184426	20.00	ng	0.00
75) Chrysene-d12	13.77	240	174024	20.00	ng	0.00
86) Perylene-d12	15.16	264	125417	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	250725	39.82	ng	0.00
7) Phenol-d6	6.27	99	359523	47.58	ng	0.00
23) Nitrobenzene-d5	7.19	82	202394	36.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.45	330	16085	15.14	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	351277	44.60	ng	0.00
78) Terphenyl-d14	12.73	244	230708	22.99	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.39	163	73091	7.800	ng	98
70) Phenanthrene	11.17	178	45417	4.529	ng	98
74) Fluoranthene	12.35	202	41166	4.289	ng	98
77) Pyrene	12.57	202	40840	2.586	ng	99
85) Indeno(1,2,3-cd)pyrene	16.47	276	6583	3.970	ng	# 88

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

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