

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050516\
 Data File : BF086718.D
 Acq On : 5 May 2016 21:16
 Operator : UM/SJ
 Sample : H2829-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EUT01-TP-B8(0-4.5)-C

Manual Integrations
 APPROVED

umangi
 5/6/2016 7:11:44 PM

Quant Time: May 06 02:27:06 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 15:12:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	194432	40.00	ng	-0.01
21) Naphthalene-d8	8.01	136	746720	40.00	ng	-0.02
38) Acenaphthene-d10	9.77	164	335487	40.00	ng	-0.01
63) Phenanthrene-d10	11.24	188	556155	40.00	ng	-0.01
75) Chrysene-d12	13.87	240	372301	40.00	ng	-0.01
86) Perylene-d12	15.27	264	364956	40.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	1123313	95.64	ng	0.00
7) Phenol-d6	6.37	99	1320091	83.44	ng	0.00
23) Nitrobenzene-d5	7.30	82	1033063	71.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.56	330	330526	100.71	ng	-0.01
44) 2-Fluorobiphenyl	9.09	172	1394124	68.12	ng	-0.01
78) Terphenyl-d14	12.83	244	1028223	55.82	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.48	163	124210	4.97	ng	99
74) Fluoranthene	12.45	202	106009	3.54	ng	92
77) Pyrene	12.67	202	101899	3.43	ng	97
80) Benzo(a)anthracene	13.86	228	57046	2.71	ng	92
82) Chrysene	13.89	228	46883m	2.09	ng	
87) Benzo(b)fluoranthene	14.88	252	77977m	3.56	ng	
89) Benzo(a)pyrene	15.21	252	49404	2.45	ng	# 92

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

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