

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086434.D
 Acq On : 27 Apr 2016 20:04
 Operator : UM/SJ
 Sample : H2724-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TH-B2(0-5)-C

Manual Integrations
 APPROVED

sohil
 4/28/2016 1:40:17 PM

Quant Time: Apr 28 01:38:45 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	121004	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	510195	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	209667	20.00	ng	-0.01
63) Phenanthrene-d10	11.32	188	338642	20.00	ng	0.00
75) Chrysene-d12	13.95	240	263243	20.00	ng	-0.01
86) Perylene-d12	15.37	264	257743	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	541816	77.23	ng	0.00
7) Phenol-d6	6.44	99	788313	80.51	ng	0.00
23) Nitrobenzene-d5	7.37	82	503012	53.14	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	132696	60.01	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	772618	60.05	ng	-0.01
78) Terphenyl-d14	12.91	244	477451	39.96	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	108124	6.85	ng	Qvalue 100
70) Phenanthrene	11.34	178	109923	6.19	ng	99
74) Fluoranthene	12.53	202	136322	7.60	ng	92
77) Pyrene	12.75	202	119568	6.30	ng	95
80) Benzo(a)anthracene	13.94	228	72846	5.19	ng	96
82) Chrysene	13.97	228	62024m	4.07	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	35644	3.15	ng	# 88
87) Benzo(b)fluoranthene	14.97	252	97470m	6.43	ng	
88) Benzo(k)fluoranthene	14.98	252	38073m	2.41	ng	
89) Benzo(a)pyrene	15.30	252	63533	4.44	ng	# 91
91) Benzo(g,h,i)perylene	17.13	276	32872	2.80	ng	95

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

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