

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086405.D
 Acq On : 23 Apr 2016 22:19
 Operator : UM/SJ
 Sample : H2652-03
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B10(5-8.5)-C

Manual Integrations
 APPROVED

sohil
 4/25/2016 12:36:11 PM

Quant Time: Apr 25 07:09:13 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 23 00:24:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	168587	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	663650	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	282260	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	469339	20.00	ng	0.00
75) Chrysene-d12	14.00	240	348714	20.00	ng	0.00
86) Perylene-d12	15.41	264	257177	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1038379	106.37	ng	0.02
7) Phenol-d6	6.49	99	1370475	103.26	ng	0.01
23) Nitrobenzene-d5	7.41	82	833407	74.39	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	237062	92.78	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1351835	81.73	ng	0.00
78) Terphenyl-d14	12.94	244	949800	67.67	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	305571	14.99	ng	99
70) Phenanthrene	11.39	178	71234	3.02	ng	99
74) Fluoranthene	12.57	202	209172	8.22	ng	95
77) Pyrene	12.80	202	190904	7.85	ng	97
80) Benzo(a)anthracene	13.99	228	81606	4.51	ng	99
82) Chrysene	14.02	228	69116	3.39	ng	98
87) Benzo(b)fluoranthene	15.01	252	75581m	5.35	ng	
88) Benzo(k)fluoranthene	15.04	252	37247m	2.69	ng	
89) Benzo(a)pyrene	15.36	252	52481	3.77	ng	# 86
91) Benzo(g,h,i)perylene	17.21	276	32652	2.75	ng	93

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

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