

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104131.D
 Acq On : 2 Apr 2018 13:29
 Operator : JU/SJ
 Sample : J2142-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GB-RT-2342

Manual Integrations
 APPROVED

Sohil
 4/3/2018 2:26:56 PM

Quant Time: Apr 02 15:47:57 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 02 13:40:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	98390	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	418299	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	198691	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	376851	20.00	ng	0.00
75) Chrysene-d12	14.16	240	266547	20.00	ng	0.01
86) Perylene-d12	15.73	264	225812	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	502474	81.44	ng	0.00
7) Phenol-d6	6.59	99	857829	113.01	ng	0.00
23) Nitrobenzene-d5	7.52	82	534460	78.14	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	178023	80.46	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	848798	67.37	ng	0.00
78) Terphenyl-d14	13.07	244	793608	68.75	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.70	163	71194	4.538	ng	100
70) Phenanthrene	11.52	178	46784	2.293	ng	95
74) Fluoranthene	12.71	202	113471	5.232	ng	98
77) Pyrene	12.94	202	95242	4.971	ng	98
80) Benzo(a)anthracene	14.15	228	49462	2.966	ng	97
82) Chrysene	14.19	228	43068	2.636	ng	97
87) Benzo(b)fluoranthene	15.27	252	43592m	3.172	ng	
89) Benzo(a)pyrene	15.66	252	32556	2.556	ng	# 90

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

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