

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041718\
 Data File : BF104566.D
 Acq On : 17 Apr 2018 17:31
 Operator : JU/SJ
 Sample : J2395-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2-1-2-DENSE-GRADE-AGG-RCARE

Manual Integrations
 APPROVED

Sohil
 4/18/2018 3:00:38 PM

Quant Time: Apr 17 18:19:51 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 17 17:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	131805	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	534251	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	230206	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	386351	20.00	ng	0.00
75) Chrysene-d12	13.97	240	314031	20.00	ng	0.00
86) Perylene-d12	15.44	264	275983	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	41582	4.82	ng	0.00
7) Phenol-d6	6.43	99	413465	39.04	ng	-0.01
23) Nitrobenzene-d5	7.37	82	609878	74.54	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	477	0.20	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1037859	71.22	ng	0.00
78) Terphenyl-d14	12.92	244	915673	66.48	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	169332	9.328	ng	98
70) Phenanthrene	11.36	178	93813	4.360	ng	97
74) Fluoranthene	12.54	202	113971	5.070	ng	98
77) Pyrene	12.77	202	106139	4.560	ng	98
80) Benzo(a)anthracene	13.96	228	49179	2.621	ng	97
82) Chrysene	14.00	228	48775	2.531	ng	97
87) Benzo(b)fluoranthene	15.02	252	56887m	3.264	ng	
89) Benzo(a)pyrene	15.37	252	37888	2.429	ng	# 86

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

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