

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041618\
 Data File : BF104541.D
 Acq On : 16 Apr 2018 19:07
 Operator : JU/SJ
 Sample : J2395-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 4/17/2018 10:09:01 AM

Quant Time: Apr 17 01:11:06 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 16 12:15:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	157907	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	642889	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	293159	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	492006	20.00	ng	0.00
75) Chrysene-d12	13.99	240	298386	20.00	ng	0.00
86) Perylene-d12	15.45	264	321887	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	50074	4.85	ng	0.00
7) Phenol-d6	6.44	99	490562	38.66	ng	-0.01
23) Nitrobenzene-d5	7.37	82	721287	73.26	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	584	0.19	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1241696	66.91	ng	0.00
78) Terphenyl-d14	12.93	244	1021453	78.04	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	215288	9.313	ng	100
70) Phenanthrene	11.36	178	122839	4.483	ng	97
74) Fluoranthene	12.55	202	133029	4.647	ng	99
77) Pyrene	12.78	202	117700	5.322	ng	97
80) Benzo(a)anthracene	13.97	228	47032	2.638	ng	95
82) Chrysene	14.01	228	46732	2.552	ng	98
87) Benzo(b)fluoranthene	15.02	252	55212m	2.716	ng	
89) Benzo(a)pyrene	15.39	252	42358	2.329	ng	# 92
91) Benzo(g,h,i)perylene	17.34	276	30732	2.123	ng	94

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

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