

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041718\
 Data File : BF104567.D
 Acq On : 17 Apr 2018 17:57
 Operator : JU/SJ
 Sample : J2393-01RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-4-DENSE-GRADE-AGG-RCARE

Manual Integrations
 APPROVED

Sohil
 4/18/2018 3:00:42 PM

Quant Time: Apr 18 01:03:02 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	132251	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	537092	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	241165	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	393021	20.00	ng	0.00
75) Chrysene-d12	13.97	240	267648	20.00	ng	0.00
86) Perylene-d12	15.44	264	270575	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	23423	2.71	ng	0.00
7) Phenol-d6	6.43	99	340414	32.03	ng	-0.01
23) Nitrobenzene-d5	7.37	82	554718	67.44	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	456	0.18	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	990246	64.87	ng	0.00
78) Terphenyl-d14	12.92	244	776537	66.15	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	121393	6.383	ng	99
70) Phenanthrene	11.36	178	133495	6.099	ng	98
74) Fluoranthene	12.55	202	150814	6.595	ng	97
77) Pyrene	12.77	202	135248	6.817	ng	98
80) Benzo(a)anthracene	13.96	228	62463	3.906	ng	98
82) Chrysene	14.00	228	54917	3.344	ng	96
85) Indeno(1,2,3-cd)pyrene	16.89	276	28304	2.029	ng	94
87) Benzo(b)fluoranthene	15.01	252	75866m	4.439	ng	
89) Benzo(a)pyrene	15.37	252	49883	3.262	ng	# 92
91) Benzo(g,h,i)perylene	17.32	276	29310	2.409	ng	# 95

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

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