

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
 Data File : BF103209.D
 Acq On : 26 Feb 2018 11:01
 Operator : SJ/JU
 Sample : J1640-01RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 2/27/2018 12:22:25 PM

Quant Time: Feb 26 12:22:07 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	150021	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	648916	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	290700	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	520044	20.00	ng	0.00
75) Chrysene-d12	14.04	240	351408	20.00	ng	0.00
86) Perylene-d12	15.53	264	289020	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	24010	2.42	ng	0.01
7) Phenol-d6	6.49	99	396175	32.64	ng	-0.01
23) Nitrobenzene-d5	7.43	82	585697	51.54	ng	0.00
41) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
44) 2-Fluorobiphenyl	9.22	172	977799	52.73	ng	0.00
78) Terphenyl-d14	12.98	244	911176	62.33	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	91857	3.939	ng	98
70) Phenanthrene	11.42	178	102557	3.583	ng	98
74) Fluoranthene	12.61	202	169700	5.901	ng	99
77) Pyrene	12.84	202	170855	6.236	ng	100
80) Benzo(a)anthracene	14.03	228	61054	2.678	ng	99
82) Chrysene	14.07	228	63890	2.886	ng	96
87) Benzo(b)fluoranthene	15.10	252	64050m	3.371	ng	
89) Benzo(a)pyrene	15.47	252	44397	2.616	ng	# 91
91) Benzo(g,h,i)perylene	17.50	276	32593	2.543	ng	95

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

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