

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060117\
 Data File : BF095627.D
 Acq On : 2 Jun 2017 9:46
 Operator : SJ/MA
 Sample : I3415-01 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-U6-SSW

Manual Integrations
 APPROVED

mohammad
 6/2/2017 2:33:09 PM

Quant Time: Jun 02 13:55:27 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 01 03:03:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	127357	20.00	ng	-0.02
21) Naphthalene-d8	9.57	136	461614	20.00	ng	-0.02
38) Acenaphthene-d10	12.39	164	213399	20.00	ng	-0.02
63) Phenanthrene-d10	14.80	188	338210	20.00	ng	-0.02
75) Chrysene-d12	18.49	240	233770	20.00	ng	-0.02
86) Perylene-d12	20.17	264	218652	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	50016m	6.55	ng	0.09
7) Phenol-d6	7.16	99	88192m	9.62	ng	0.05
23) Nitrobenzene-d5	8.50	82	38963	5.32	ng	0.03
41) 2,4,6-Tribromophenol	13.73	330	6594	3.77	ng	0.02
44) 2-Fluorobiphenyl	11.35	172	102823	6.94	ng	-0.02
78) Terphenyl-d14	17.14	244	59956	5.65	ng	-0.02
Target Compounds						
49) Dimethylphthalate	12.06	163	36845	2.103	ng	Qvalue 97
70) Phenanthrene	14.84	178	55689	2.866	ng	97
74) Fluoranthene	16.60	202	86110	4.320	ng	97
77) Pyrene	16.91	202	77762	4.448	ng	95
80) Benzo(a)anthracene	18.49	228	29176	2.144	ng	98
82) Chrysene	18.53	228	29713	2.259	ng	94
87) Benzo(b)fluoranthene	19.77	252	35099	2.598	ng	# 96
89) Benzo(a)pyrene	20.12	252	27171	2.179	ng	# 93

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

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