

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060117\
 Data File : BF095628.D
 Acq On : 2 Jun 2017 10:18
 Operator : SJ/MA
 Sample : I3415-03 10X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-T6-ESW

Manual Integrations
 APPROVED

mohammad
 6/2/2017 2:33:11 PM

Quant Time: Jun 02 13:58:27 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	111794	20.00	ng	-0.02
21) Naphthalene-d8	9.57	136	434225	20.00	ng	-0.02
38) Acenaphthene-d10	12.39	164	188335	20.00	ng	-0.02
63) Phenanthrene-d10	14.80	188	291828	20.00	ng	-0.02
75) Chrysene-d12	18.49	240	221041	20.00	ng	-0.02
86) Perylene-d12	20.17	264	186577	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	61262m	9.14	ng	0.07
7) Phenol-d6	7.16	99	85456m	10.62	ng	0.05
23) Nitrobenzene-d5	8.50	82	44716	6.49	ng	0.03
41) 2,4,6-Tribromophenol	13.72	330	11381	7.38	ng	0.01
44) 2-Fluorobiphenyl	11.35	172	97414	7.44	ng	-0.02
78) Terphenyl-d14	17.14	244	55826	5.56	ng	-0.02
Target Compounds						
70) Phenanthrene	14.84	178	83735	4.994	ng	99
74) Fluoranthene	16.60	202	107551	6.253	ng	98
77) Pyrene	16.91	202	89804	5.432	ng	# 95
80) Benzo(a)anthracene	18.49	228	34564	2.687	ng	97
82) Chrysene	18.53	228	33761	2.714	ng	100
87) Benzo(b)fluoranthene	19.77	252	37239	3.230	ng	# 96
89) Benzo(a)pyrene	20.12	252	28102	2.642	ng	# 95

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

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