

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051517\
 Data File : BF095137.D
 Acq On : 15 May 2017 21:23
 Operator : SJ/MA
 Sample : I3147-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3COMP

Manual Integrations
 APPROVED

mohammad
 5/17/2017 2:20:12 PM

Quant Time: May 16 08:21:05 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	95648	20.00	ng	0.05
21) Naphthalene-d8	9.80	136	380351	20.00	ng	0.04
38) Acenaphthene-d10	12.62	164	167407	20.00	ng	0.04
63) Phenanthrene-d10	15.03	188	259554	20.00	ng	0.05
75) Chrysene-d12	18.70	240	200545	20.00	ng	0.05
86) Perylene-d12	20.38	264	166632	20.00	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	799657	139.39	ng	0.07
7) Phenol-d6	7.32	99	1037173	150.67	ng	0.05
23) Nitrobenzene-d5	8.68	82	596119	98.83	ng	0.04
41) 2,4,6-Tribromophenol	13.92	330	178986	130.55	ng	0.04
44) 2-Fluorobiphenyl	11.57	172	1150773	98.94	ng	0.04
78) Terphenyl-d14	17.35	244	738136	81.07	ng	0.04
Target Compounds						Qvalue
49) Dimethylphthalate	12.26	163	170243	12.39	ng	98
77) Pyrene	17.12	202	34123	2.28	ng	95
80) Benzo(a)anthracene	18.69	228	24519	2.10	ng	# 91
82) Chrysene	18.74	228	28450	2.52	ng	# 94
85) Indeno(1,2,3-cd)pyrene	21.73	276	20823	2.27	ng	91
87) Benzo(b)fluoranthene	19.97	252	38258m	3.72	ng	
89) Benzo(a)pyrene	20.32	252	32703	3.44	ng	# 89
91) Benzo(g,h,i)perylene	22.12	276	23877	3.10	ng	90

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

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