

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071317\
 Data File : BF096754.D
 Acq On : 14 Jul 2017 3:41
 Operator : SJ/JU
 Sample : I4172-01
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-P6-SSW

Manual Integrations
 APPROVED

mohammad
 7/17/2017 8:12:04 AM

Quant Time: Jul 14 06:24:25 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 14:49:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.64	152	114851	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	422492	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	140714	20.00	ng	0.00
63) Phenanthrene-d10	11.15	188	246897	20.00	ng	0.00
75) Chrysene-d12	13.77	240	196505	20.00	ng	0.00
86) Perylene-d12	15.16	264	135516	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	766770	100.38	ng	0.02
7) Phenol-d6	6.29	99	895516	97.70	ng	0.00
23) Nitrobenzene-d5	7.20	82	486503	71.33	ng	0.00
41) 2,4,6-Tribromophenol	10.46	330	127724	106.34	ng	0.00
44) 2-Fluorobiphenyl	9.00	172	670406	75.27	ng	0.00
78) Terphenyl-d14	12.73	244	447450	39.49	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.39	163	88342	8.338	ng	97
70) Phenanthrene	11.17	178	57555	4.287	ng	98
74) Fluoranthene	12.35	202	68798	5.354	ng	97
77) Pyrene	12.58	202	63176	3.542	ng	98
80) Benzo(a)anthracene	13.76	228	31339	2.560	ng	96
82) Chrysene	13.80	228	27340	2.268	ng	96
87) Benzo(b)fluoranthene	14.77	252	24989m	3.057	ng	
89) Benzo(a)pyrene	15.10	252	18274	2.438	ng	# 93

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

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