

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070717\
 Data File : BF096576.D
 Acq On : 7 Jul 2017 23:12
 Operator : SJ/MA
 Sample : I4070-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-070517-C

Manual Integrations
 APPROVED

mohammad
 7/10/2017 2:22:51 PM

Quant Time: Jul 10 03:16:23 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 13:02:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	114660	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	456920	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	189401	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	259645	20.00	ng	0.00
75) Chrysene-d12	13.83	240	213885	20.00	ng	-0.01
86) Perylene-d12	15.23	264	221540	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	744070	95.78	ng	0.00
7) Phenol-d6	6.34	99	881503	92.30	ng	0.00
23) Nitrobenzene-d5	7.26	82	512615	67.42	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	143388	96.92	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	835034	75.38	ng	0.00
78) Terphenyl-d14	12.79	244	477650	47.72	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.35	94	27273	2.507	ng	93
49) Dimethylphthalate	9.45	163	98182	6.944	ng	99
70) Phenanthrene	11.23	178	71530	5.096	ng	97
74) Fluoranthene	12.41	202	62926	4.573	ng	96
77) Pyrene	12.64	202	86928	5.063	ng	97
80) Benzo(a)anthracene	13.82	228	31953m	2.580	ng	
82) Chrysene	13.86	228	42599	3.284	ng	97
87) Benzo(b)fluoranthene	14.84	252	41796m	3.011	ng	
89) Benzo(a)pyrene	15.17	252	32604	2.631	ng	92

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

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