

Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
 Data File : BF095583.D
 Acq On : 1 Jun 2017 1:29
 Operator : SJ/MA
 Sample : I3302-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM15-COMP-6-18

Manual Integrations
 APPROVED

mohammad
 6/1/2017 2:20:46 PM

Quant Time: Jun 01 07:35:09 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 04:09:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	89826	20.00	ng	0.00
21) Naphthalene-d8	9.58	136	393393	20.00	ng	0.00
38) Acenaphthene-d10	12.41	164	138193	20.00	ng	0.00
63) Phenanthrene-d10	14.81	188	208834	20.00	ng	0.00
75) Chrysene-d12	18.51	240	175574	20.00	ng	0.00
86) Perylene-d12	20.18	264	133647	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	597045	110.81	ng	0.00
7) Phenol-d6	7.11	99	751954	116.32	ng	0.00
23) Nitrobenzene-d5	8.47	82	422647	67.75	ng	0.00
41) 2,4,6-Tribromophenol	13.71	330	100707	88.98	ng	0.00
44) 2-Fluorobiphenyl	11.36	172	598110	62.30	ng	0.00
78) Terphenyl-d14	17.15	244	355214	44.56	ng	0.00
Target Compounds						
49) Dimethylphthalate	12.05	163	103235	9.10	ng	Qvalue # 98
70) Phenanthrene	14.85	178	65888	5.49	ng	96
74) Fluoranthene	16.61	202	112845	9.17	ng	99
77) Pyrene	16.91	202	103598	7.89	ng	95
80) Benzo(a)anthracene	18.49	228	52651	5.15	ng	97
82) Chrysene	18.54	228	48980	4.96	ng	98
85) Indeno(1,2,3-cd)pyrene	21.44	276	21738	2.71	ng	99
87) Benzo(b)fluoranthene	19.78	252	56517m	6.84	ng	
88) Benzo(k)fluoranthene	19.81	252	19985m	2.51	ng	
89) Benzo(a)pyrene	20.13	252	40184	5.27	ng	98
91) Benzo(g,h,i)perylene	21.80	276	20536	3.33	ng	92

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

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