

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052617\
 Data File : BF095490.D
 Acq On : 27 May 2017 8:11
 Operator : SJ/MA
 Sample : I3339-03 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-S10-BASE

Manual Integrations
 APPROVED

mohammad
 5/30/2017 1:37:25 PM

Quant Time: May 28 23:33:58 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 26 16:50:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	82137	20.00	ng	0.00
21) Naphthalene-d8	9.77	136	323844	20.00	ng	0.00
38) Acenaphthene-d10	12.59	164	133197	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	196755	20.00	ng	0.00
75) Chrysene-d12	18.68	240	160662	20.00	ng	0.00
86) Perylene-d12	20.36	264	136668	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.77	112	104313	21.17	ng	0.06
7) Phenol-d6	7.34	99	119888	20.28	ng	0.05
23) Nitrobenzene-d5	8.68	82	73966	14.40	ng	0.02
41) 2,4,6-Tribromophenol	13.92	330	17571	16.11	ng	0.02
44) 2-Fluorobiphenyl	11.54	172	131284	14.19	ng	0.00
78) Terphenyl-d14	17.32	244	73208	10.04	ng	0.00
Target Compounds						
31) Naphthalene	9.80	128	80773	4.96	ng	97
49) Dimethylphthalate	12.24	163	60693	5.55	ng	96
70) Phenanthrene	15.05	178	41961	3.71	ng	98
74) Fluoranthene	16.79	202	40548	3.50	ng	96
77) Pyrene	17.09	202	63510	5.29	ng	96
80) Benzo(a)anthracene	18.67	228	19308	2.06	ng	98
82) Chrysene	18.71	228	23486m	2.60	ng	
87) Benzo(b)fluoranthene	19.96	252	24994m	2.96	ng	
89) Benzo(a)pyrene	20.31	252	24649	3.16	ng	97
91) Benzo(g,h,i)perylene	22.11	276	17136	2.71	ng	94

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

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