

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
 Data File : BF114552.D
 Acq On : 24 May 2019 4:12
 Operator : JU/SJ
 Sample : K3008-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-21

Manual Integrations
 APPROVED

mohammad
 5/24/2019 1:36:59 PM

Quant Time: May 24 08:33:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 24 07:23:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	158532	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	642003	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	314586	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	593591	20.00	ng	0.00
76) Chrysene-d12	13.98	240	358620	20.00	ng	0.00
87) Perylene-d12	15.44	264	333389	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	409678	41.59	ng	0.00
7) Phenol-d6	6.46	99	595846	51.14	ng	0.00
23) Nitrobenzene-d5	7.37	82	359088	31.39	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	109994	33.82	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	706255	33.96	ng	0.00
79) Terphenyl-d14	12.92	244	625589	35.96	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	104952	4.545	ng	98
71) Phenanthrene	11.36	178	263087	9.110	ng	97
75) Fluoranthene	12.55	202	375094	12.061	ng	100
78) Pyrene	12.78	202	327396	11.349	ng	99
81) Benzo(a)anthracene	13.97	228	155860	6.719	ng	99
83) Chrysene	14.01	228	146971	6.464	ng	96
86) Indeno(1,2,3-cd)pyrene	16.92	276	63926	3.181	ng	# 94
88) Benzo(b)fluoranthene	15.02	252	178299m	8.348	ng	
89) Benzo(k)fluoranthene	15.04	252	60907m	3.104	ng	
90) Benzo(a)pyrene	15.39	252	120464	6.409	ng	# 94
92) Benzo(g,h,i)perylene	17.37	276	60061	3.837	ng	97

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

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