

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114961.D
 Acq On : 11 Jun 2019 23:20
 Operator : HP/JU
 Sample : K3247-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-041-A

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:09:07 AM

Quant Time: Jun 12 05:32:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 16:44:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	238626	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	918618	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	448190	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	835822	20.00	ng	0.00
76) Chrysene-d12	13.79	240	615228	20.00	ng	0.00
87) Perylene-d12	15.17	264	556195	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1241627	88.06	ng	0.00
7) Phenol-d6	6.30	99	1533820	90.19	ng	0.00
23) Nitrobenzene-d5	7.22	82	907536	60.70	ng	0.00
42) 2,4,6-Tribromophenol	10.47	330	404418	82.95	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	1770866	62.71	ng	0.00
79) Terphenyl-d14	12.74	244	1786600	57.75	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.41	163	227833	6.763	ng	98
71) Phenanthrene	11.19	178	253372	6.248	ng	95
75) Fluoranthene	12.36	202	404937	9.654	ng	96
78) Pyrene	12.59	202	374655	6.992	ng	97
81) Benzo(a)anthracene	13.77	228	193072	4.244	ng	97
83) Chrysene	13.81	228	175484	3.841	ng	97
88) Benzo(b)fluoranthene	14.78	252	214915m	5.935	ng	
89) Benzo(k)fluoranthene	14.80	252	67880m	2.160	ng	
90) Benzo(a)pyrene	15.11	252	134514	4.277	ng	# 94
92) Benzo(g,h,i)perylene	16.86	276	70046	2.645	ng	# 91

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

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