

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061820\
 Data File : BF120645.D
 Acq On : 18 Jun 2020 16:43
 Operator : JU/CG
 Sample : L3046-11
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ISS-B-C

Manual Integrations
 APPROVED

mohammad
 6/19/2020 12:57:54 PM

Quant Time: Jun 18 18:36:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	115082	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	412171	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	263928	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	455228	20.00	ng	0.00
76) Chrysene-d12	13.96	240	362253	20.00	ng	0.00
86) Perylene-d12	15.40	264	442010	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	528886	78.79	ng	0.00
7) Phenol-d6	6.43	99	788387	94.30	ng	0.00
23) Nitrobenzene-d5	7.36	82	435995	59.13	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	205201	66.87	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	889659	51.57	ng	0.00
79) Terphenyl-d14	12.90	244	1056356	64.41	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	90007	4.735	ng	99
71) Phenanthrene	11.34	178	131364	5.805	ng	98
75) Fluoranthene	12.53	202	204602	7.908	ng	97
78) Pyrene	12.76	202	214699	8.366	ng	98
81) Benzo(a)anthracene	13.94	228	81870	3.762	ng	98
83) Chrysene	13.98	228	90564	4.057	ng	96
87) Indeno(1,2,3-cd)pyrene	16.81	276	54145	2.014	ng	97
88) Benzo(b)fluoranthene	14.98	252	122513m	4.653	ng	
90) Benzo(a)pyrene	15.33	252	76477	3.246	ng	# 92
92) Benzo(g,h,i)perylene	17.24	276	48023	2.210	ng	98

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

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