

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
 Data File : BF120561.D
 Acq On : 5 Jun 2020 15:58
 Operator : JU/CG
 Sample : L2890-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11979-72-11969

Manual Integrations
 APPROVED

mohammad
 6/8/2020 12:51:07 PM

Quant Time: Jun 05 17:34:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:36:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	128281	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	487118	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	228627	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	347464	20.00	ng	0.00
76) Chrysene-d12	13.98	240	275250	20.00	ng	0.00
86) Perylene-d12	15.43	264	271907	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	973063	144.56	ng	0.00
7) Phenol-d6	6.46	99	742996	89.55	ng	0.00
23) Nitrobenzene-d5	7.38	82	509684	68.44	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	193718	91.45	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	984362	73.05	ng	0.00
79) Terphenyl-d14	12.93	244	835771	67.57	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.57	163	187761	13.136	ng	99
71) Phenanthrene	11.36	178	58400	3.814	ng	# 89
75) Fluoranthene	12.55	202	104211	6.326	ng	97
78) Pyrene	12.78	202	94721	4.948	ng	97
81) Benzo(a)anthracene	13.97	228	48989	3.321	ng	98
83) Chrysene	14.00	228	53643m	3.469	ng	
87) Indeno(1,2,3-cd)pyrene	16.86	276	34108	2.293	ng	97
88) Benzo(b)fluoranthene	15.00	252	65920m	4.365	ng	
90) Benzo(a)pyrene	15.36	252	41752	3.161	ng	# 78
92) Benzo(g,h,i)perylene	17.29	276	34454	2.881	ng	# 88

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

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