

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
 Data File : BF119880.D
 Acq On : 9 Apr 2020 20:05
 Operator : MA/SJ
 Sample : L2166-05
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-47-COMP

Manual Integrations
 APPROVED

mohammad
 4/10/2020 9:57:49 AM

Quant Time: Apr 10 02:39:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 09 01:21:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	159755	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	625725	20.00	ng	0.00
39) Acenaphthene-d10	9.81	164	272350	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	408924	20.00	ng	0.00
76) Chrysene-d12	13.93	240	287147	20.00	ng	0.00
87) Perylene-d12	15.37	264	290564	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1397056	151.13	ng	0.01
7) Phenol-d6	6.41	99	1815869	152.45	ng	0.00
23) Nitrobenzene-d5	7.33	82	1096341	90.64	ng	0.00
42) 2,4,6-Tribromophenol	10.60	330	373407	111.98	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	1818973	94.82	ng	0.00
79) Terphenyl-d14	12.89	244	1451328	85.05	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.53	163	331856	15.753	ng	98
75) Fluoranthene	12.51	202	59847	2.549	ng	96
78) Pyrene	12.73	202	114135	4.715	ng	97
81) Benzo(a)anthracene	13.93	228	104606	5.538	ng	96
83) Chrysene	13.96	228	110771	5.771	ng	99
86) Indeno(1,2,3-cd)pyrene	16.79	276	53792	3.179	ng	96
88) Benzo(b)fluoranthene	14.96	252	120913m	5.785	ng	
90) Benzo(a)pyrene	15.32	252	113729	6.471	ng	98
92) Benzo(g,h,i)perylene	17.21	276	56407	3.671	ng	99

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

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