

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100520\
 Data File : BF121834.D
 Acq On : 5 Oct 2020 23:58
 Operator : JU/CG
 Sample : L4226-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-100120

Manual Integrations
 APPROVED

mohammad
 10/7/2020 10:36:37 AM

Quant Time: Oct 06 02:35:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	107195	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	396806	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	182574	20.00	ng	-0.01
64) Phenanthrene-d10	11.28	188	299023	20.00	ng	-0.01
76) Chrysene-d12	13.92	240	284899	20.00	ng	-0.01
86) Perylene-d12	15.36	264	265533	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	271958	37.70	ng	0.00
7) Phenol-d6	6.40	99	341485	36.09	ng	-0.01
23) Nitrobenzene-d5	7.33	82	221203	29.93	ng	-0.01
42) 2,4,6-Tribromophenol	10.59	330	71765	31.63	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	385416	31.97	ng	-0.01
79) Terphenyl-d14	12.86	244	315581	22.44	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.52	163	50535	3.819	ng	99
75) Fluoranthene	12.50	202	99807	5.739	ng	96
78) Pyrene	12.72	202	90190	4.333	ng	98
81) Benzo(a)anthracene	13.91	228	50633	2.809	ng	96
83) Chrysene	13.95	228	52221	2.861	ng	97
88) Benzo(b)fluoranthene	14.94	252	67151m	3.783	ng	
90) Benzo(a)pyrene	15.30	252	43094	2.856	ng	# 92

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

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