

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
 Data File : BF119888.D
 Acq On : 9 Apr 2020 23:42
 Operator : MA/SJ
 Sample : L2157-01
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-226

Manual Integrations
 APPROVED

mohammad
 4/10/2020 9:57:59 AM

Quant Time: Apr 10 04:06:45 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	170361	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	581666	20.00	ng	0.00
39) Acenaphthene-d10	9.81	164	235375	20.00	ng	0.00
64) Phenanthrene-d10	11.30	188	359746	20.00	ng	0.00
76) Chrysene-d12	13.94	240	286165	20.00	ng	0.00
87) Perylene-d12	15.37	264	250658	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1473577	149.48	ng	0.01
7) Phenol-d6	6.41	99	1821263	143.38	ng	0.00
23) Nitrobenzene-d5	7.33	82	1133425	100.81	ng	0.00
42) 2,4,6-Tribromophenol	10.60	330	333572	115.75	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	1522054	91.81	ng	0.00
79) Terphenyl-d14	12.89	244	1221374	71.82	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.53	163	325654	17.887	ng	98
71) Phenanthrene	11.32	178	49315	2.576	ng	98
75) Fluoranthene	12.51	202	109781	5.314	ng	95
78) Pyrene	12.74	202	98964	4.102	ng	99
81) Benzo(a)anthracene	13.93	228	42864	2.277	ng	94
83) Chrysene	13.96	228	47358	2.476	ng	99
88) Benzo(b)fluoranthene	14.96	252	60649m	3.363	ng	
90) Benzo(a)pyrene	15.32	252	36054	2.378	ng	# 91

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

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