

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100220\
 Data File : BF121796.D
 Acq On : 2 Oct 2020 16:11
 Operator : JU/CG
 Sample : L4198-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-25

Manual Integrations
 APPROVED

mohammad
 10/5/2020 10:27:26 AM

Quant Time: Oct 02 17:30:47 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	145775	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	570064	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	274084	20.00	ng	-0.01
64) Phenanthrene-d10	11.28	188	456449	20.00	ng	-0.01
76) Chrysene-d12	13.92	240	286097	20.00	ng	-0.02
86) Perylene-d12	15.35	264	280408	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	747746	76.23	ng	0.00
7) Phenol-d6	6.40	99	940098	73.05	ng	-0.01
23) Nitrobenzene-d5	7.33	82	649230	61.15	ng	-0.01
42) 2,4,6-Tribromophenol	10.59	330	242792	71.27	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1088610	60.15	ng	-0.01
79) Terphenyl-d14	12.86	244	849888	60.18	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.52	163	105659	5.319	ng	98
75) Fluoranthene	12.49	202	108976	4.105	ng	99
78) Pyrene	12.72	202	92084	4.405	ng	98
81) Benzo(a)anthracene	13.90	228	39745	2.196	ng	98
83) Chrysene	13.94	228	38315	2.090	ng	98
88) Benzo(b)fluoranthene	14.94	252	46648m	2.488	ng	

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

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