

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\
 Data File : BF121483.D
 Acq On : 31 Aug 2020 18:00
 Operator : JU/CG
 Sample : L3508-25
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-08-082820

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:56:04 PM

Quant Time: Aug 31 18:41:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	326533	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1189794	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	590221	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	972364	20.00	ng	0.00
76) Chrysene-d12	14.02	240	678160	20.00	ng	0.00
86) Perylene-d12	15.49	264	721205	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1506353	77.40	ng	0.00
7) Phenol-d6	6.48	99	1971813	72.50	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1276067	73.32	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	501196	84.07	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2241429	61.28	ng	0.00
79) Terphenyl-d14	12.97	244	1801808	55.93	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	179685	4.432	ng	99
75) Fluoranthene	12.59	202	247504	4.602	ng	98
78) Pyrene	12.82	202	300144	6.197	ng	96
81) Benzo(a)anthracene	14.01	228	206771	4.995	ng	99
83) Chrysene	14.04	228	186181	4.495	ng	95
87) Indeno(1,2,3-cd)pyrene	16.95	276	95662	2.156	ng	98
88) Benzo(b)fluoranthene	15.06	252	279054m	5.948	ng	
90) Benzo(a)pyrene	15.43	252	179731	4.261	ng	# 96
92) Benzo(g,h,i)perylene	17.39	276	99252	2.837	ng	96

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

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