

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082820\
 Data File : BF121467.D
 Acq On : 28 Aug 2020 19:01
 Operator : JU/CG
 Sample : L3837-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 789UMR-TP05-1218-01

Manual Integrations
 APPROVED

mohammad
 8/31/2020 10:27:31 AM

Quant Time: Sep 04 15:58:53 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	312069	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1229932	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	641356	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	1208182	20.00	ng	0.00
76) Chrysene-d12	14.02	240	861648	20.00	ng	0.00
86) Perylene-d12	15.49	264	779950	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1408914	75.75	ng	0.00
7) Phenol-d6	6.48	99	1879760	72.32	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1225391	68.11	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	544322	84.03	ng	-0.01
45) 2-Fluorobiphenyl	9.22	172	2253336	56.69	ng	0.00
79) Terphenyl-d14	12.97	244	2318283	56.63	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	143841	3.265	ng	98
71) Phenanthrene	11.40	178	207228	3.342	ng	97
75) Fluoranthene	12.59	202	494753	7.404	ng	96
78) Pyrene	12.82	202	473959	7.702	ng	96
81) Benzo(a)anthracene	14.01	228	285798	5.433	ng	95
83) Chrysene	14.04	228	228317	4.339	ng	96
88) Benzo(b)fluoranthene	15.06	252	264088m	5.205	ng	
90) Benzo(a)pyrene	15.42	252	190079	4.166	ng	99
92) Benzo(g,h,i)perylene	17.38	276	111629	2.950	ng	97

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

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