

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\
 Data File : BF121497.D
 Acq On : 1 Sep 2020 15:29
 Operator : JU/CG
 Sample : L3848-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-BH-01-A

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:57:11 PM

Quant Time: Sep 01 17:28:03 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	6.85	152	259095	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	983754	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	516472	20.00	ng	-0.01
64) Phenanthrene-d10	11.38	188	978769	20.00	ng	0.00
76) Chrysene-d12	14.02	240	742147	20.00	ng	0.00
86) Perylene-d12	15.50	264	737675	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	2262	0.15	ng	-0.01
7) Phenol-d6	6.47	99	27788	1.29	ng	-0.02
23) Nitrobenzene-d5	7.41	82	752717	52.31	ng	-0.02
42) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
45) 2-Fluorobiphenyl	9.22	172	1469859	45.92	ng	-0.01
79) Terphenyl-d14	12.97	244	1468585	41.65	ng	0.00
Target Compounds						
25) Isophorone	7.67	82	81402	2.515	ng	98
50) Dimethylphthalate	9.61	163	77678	2.190	ng	97
75) Fluoranthene	12.59	202	235541	4.351	ng	96
78) Pyrene	12.82	202	244026	4.604	ng	96
81) Benzo(a)anthracene	14.01	228	134876	2.977	ng	96
83) Chrysene	14.04	228	122514	2.703	ng	96
88) Benzo(b)fluoranthene	15.06	252	162546m	3.387	ng	
90) Benzo(a)pyrene	15.43	252	119861	2.778	ng	97
92) Benzo(g,h,i)perylene	17.39	276	79556	2.223	ng	96

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

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