

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082820\
 Data File : BF121463.D
 Acq On : 28 Aug 2020 16:56
 Operator : JU/CG
 Sample : L3837-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 789UMR-TP11-0612-01

Manual Integrations
 APPROVED

mohammad
 8/31/2020 10:27:24 AM

Quant Time: Aug 28 17:55:27 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	310051	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1206583	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	626804	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	1165226	20.00	ng	0.00
76) Chrysene-d12	14.02	240	830627	20.00	ng	0.00
86) Perylene-d12	15.49	264	738292	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1492759	80.78	ng	0.00
7) Phenol-d6	6.48	99	1937780	75.03	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1228702	69.62	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	549752	86.83	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2225830	57.30	ng	0.00
79) Terphenyl-d14	12.97	244	2235269	56.65	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	241715	5.614	ng	98
71) Phenanthrene	11.40	178	322961	5.401	ng	96
75) Fluoranthene	12.59	202	432984	6.719	ng	96
78) Pyrene	12.82	202	343337	5.788	ng	96
81) Benzo(a)anthracene	14.01	228	173688	3.425	ng	98
83) Chrysene	14.05	228	166202	3.276	ng	94
88) Benzo(b)fluoranthene	15.06	252	197280m	4.108	ng	
90) Benzo(a)pyrene	15.42	252	131246	3.039	ng	100
92) Benzo(g,h,i)perylene	17.37	276	84037	2.346	ng	97

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

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