

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072420\
 Data File : BF121016.D
 Acq On : 24 Jul 2020 21:51
 Operator : JU/CG
 Sample : L3382-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CTR-023

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:19:49 PM

Quant Time: Jul 25 01:18:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	149026	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	513107	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	254649	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	441820	20.00	ng	0.00
76) Chrysene-d12	13.97	240	364248	20.00	ng	0.00
86) Perylene-d12	15.42	264	402860	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	799856	87.99	ng	0.00
7) Phenol-d6	6.45	99	952271	81.10	ng	0.00
23) Nitrobenzene-d5	7.37	82	591048	66.65	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	242464	86.99	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1153667	67.64	ng	0.00
79) Terphenyl-d14	12.92	244	948407	56.60	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	225804	12.291	ng	100
71) Phenanthrene	11.36	178	118746	5.226	ng	97
75) Fluoranthene	12.55	202	192979	7.662	ng	96
78) Pyrene	12.77	202	167130	6.490	ng	98
81) Benzo(a)anthracene	13.96	228	80062m	3.630	ng	
83) Chrysene	13.99	228	101089m	4.361	ng	
88) Benzo(b)fluoranthene	14.99	252	141399m	5.807	ng	
90) Benzo(a)pyrene	15.35	252	82691	3.722	ng	97
92) Benzo(g,h,i)perylene	17.26	276	47222	2.093	ng	95

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

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