

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062220\
 Data File : BF120687.D
 Acq On : 22 Jun 2020 20:35
 Operator : JU/CG
 Sample : L3082-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-14

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:03:42 PM

Quant Time: Jun 23 02:17:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	122121	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	449409	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	241097	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	453101	20.00	ng	0.00
76) Chrysene-d12	13.96	240	435949	20.00	ng	0.00
86) Perylene-d12	15.40	264	463281	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	774490	108.73	ng	-0.01
7) Phenol-d6	6.43	99	866623	97.68	ng	0.00
23) Nitrobenzene-d5	7.36	82	516649	64.26	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	335955	119.85	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1293194	82.07	ng	0.00
79) Terphenyl-d14	12.91	244	1614693	81.81	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	90868	5.233	ng	100
71) Phenanthrene	11.34	178	54815	2.434	ng	98
75) Fluoranthene	12.53	202	162634	6.315	ng	95
78) Pyrene	12.76	202	179221	5.803	ng	97
81) Benzo(a)anthracene	13.95	228	79208	3.024	ng	96
83) Chrysene	13.99	228	86785	3.230	ng	98
88) Benzo(b)fluoranthene	14.99	252	119655m	4.336	ng	
90) Benzo(a)pyrene	15.34	252	74681	3.024	ng	100
92) Benzo(g,h,i)perylene	17.25	276	50507	2.217	ng	98

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

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