

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
 Data File : BF121106.D
 Acq On : 29 Jul 2020 17:23
 Operator : JU/CG
 Sample : L3185-01 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-05-072720

Manual Integrations
 APPROVED

mohammad
 7/30/2020 3:33:34 PM

Quant Time: Jul 29 18:18:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 14:55:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	160842	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	601671	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	304356	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	539004	20.00	ng	0.00
76) Chrysene-d12	13.96	240	361632	20.00	ng	0.00
86) Perylene-d12	15.40	264	422485	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	390818	39.83	ng	0.00
7) Phenol-d6	6.44	99	470225	37.10	ng	0.00
23) Nitrobenzene-d5	7.36	82	299602	28.81	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	127268	38.20	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	588708	28.88	ng	0.00
79) Terphenyl-d14	12.90	244	489437	29.42	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.55	163	90109	4.104	ng	99
71) Phenanthrene	11.34	178	560740	20.228	ng	99
72) Anthracene	11.39	178	140767	5.057	ng	97
75) Fluoranthene	12.53	202	867785	28.242	ng	98
78) Pyrene	12.76	202	701110	27.422	ng	98
81) Benzo(a)anthracene	13.94	228	376452	17.191	ng	99
83) Chrysene	13.98	228	330587	14.365	ng	96
87) Indeno(1,2,3-cd)pyrene	16.82	276	221340	7.533	ng	95
88) Benzo(b)fluoranthene	14.99	252	450963m	17.659	ng	
89) Benzo(k)fluoranthene	15.01	252	136635m	5.867	ng	
90) Benzo(a)pyrene	15.34	252	311513	13.370	ng	98
91) Dibenzo(a,h)anthracene	16.83	278	56503m	2.354	ng	
92) Benzo(a,h,i)perylene	17.25	276	185705	7.847	ng	99

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

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