

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090820\
 Data File : BF121540.D
 Acq On : 9 Sep 2020 00:37
 Operator : JU/CG
 Sample : L3826-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BKG-3-(0.0-0.5)

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:49:43 PM

Quant Time: Sep 09 07:18:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 17:52:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	336271	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1266814	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	650406	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	1160676	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	782201	20.00	ng	-0.02
86) Perylene-d12	15.46	264	778467	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1494098	74.54	ng	0.00
7) Phenol-d6	6.47	99	1925038	68.73	ng	-0.01
23) Nitrobenzene-d5	7.40	82	1263847	68.21	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	582041	88.60	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	2233254	55.40	ng	-0.01
79) Terphenyl-d14	12.94	244	1908810	51.37	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.59	163	152269	3.408	ng	99
71) Phenanthrene	11.38	178	215063	3.611	ng	98
75) Fluoranthene	12.57	202	439084	6.840	ng	98
78) Pyrene	12.80	202	403490	7.223	ng	96
81) Benzo(a)anthracene	13.99	228	161357	3.379	ng	95
83) Chrysene	14.02	228	192060	4.020	ng	96
87) Indeno(1,2,3-cd)pyrene	16.89	276	112316	2.345	ng	93
88) Benzo(b)fluoranthene	15.03	252	229367m	4.529	ng	
90) Benzo(a)pyrene	15.39	252	154167	3.386	ng	98
92) Benzo(g,h,i)perylene	17.33	276	109239	2.892	ng	98

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

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