

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120486.D
 Acq On : 2 Jun 2020 11:18
 Operator : JU/CG
 Sample : L2792-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM107-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:35:41 PM

Quant Time: Jun 02 11:51:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 11:37:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	191043	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	698523	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	327984	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	549099	20.00	ng	0.00
76) Chrysene-d12	14.01	240	410340	20.00	ng	0.00
86) Perylene-d12	15.47	264	458880	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	851846	84.98	ng	0.00
7) Phenol-d6	6.47	99	1055017	85.38	ng	0.00
23) Nitrobenzene-d5	7.40	82	682838	63.94	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	258701	85.13	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	1251053	64.72	ng	0.00
79) Terphenyl-d14	12.95	244	1125488	61.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	224107	10.930	ng	99
71) Phenanthrene	11.39	178	314947	13.017	ng	98
72) Anthracene	11.44	178	84134	3.466	ng	98
75) Fluoranthene	12.58	202	477291	18.333	ng	98
78) Pyrene	12.81	202	401289	14.062	ng	98
81) Benzo(a)anthracene	14.00	228	219781	9.994	ng	98
83) Chrysene	14.03	228	200256	8.686	ng	96
87) Indeno(1,2,3-cd)pyrene	16.92	276	116363	4.636	ng	95
88) Benzo(b)fluoranthene	15.04	252	281782m	11.056	ng	
89) Benzo(k)fluoranthene	15.07	252	69769m	2.932	ng	
90) Benzo(a)pyrene	15.40	252	186597	8.370	ng	99
92) Benzo(a,h,i)perylene	17.36	276	107621	5.333	ng	99

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

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