

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120498.D
 Acq On : 2 Jun 2020 16:45
 Operator : JU/CG
 Sample : L2798-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM105-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:36:08 PM

Quant Time: Jun 02 17:37:02 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	162031	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	546827	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	244300	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	444864	20.00	ng	0.00
76) Chrysene-d12	14.01	240	404706	20.00	ng	0.00
86) Perylene-d12	15.47	264	332891	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	787195	92.59	ng	0.00
7) Phenol-d6	6.47	99	934762	89.20	ng	0.00
23) Nitrobenzene-d5	7.40	82	559187	66.89	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	214556	94.79	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	1015323	70.52	ng	0.00
79) Terphenyl-d14	12.96	244	1102196	60.60	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	200096	13.101	ng	Qvalue 100
71) Phenanthrene	11.39	178	47490	2.423	ng	99
75) Fluoranthene	12.58	202	116225	5.510	ng	97
78) Pyrene	12.81	202	118606	4.214	ng	97
81) Benzo(a)anthracene	14.00	228	70087	3.231	ng	95
83) Chrysene	14.03	228	59965	2.637	ng	95
88) Benzo(b)fluoranthene	15.04	252	71408m	3.862	ng	
90) Benzo(a)pyrene	15.40	252	46958	2.903	ng	# 94

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

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