

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081120\
 Data File : BF121240.D
 Acq On : 11 Aug 2020 15:14
 Operator : JU/CG
 Sample : L3623-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STP-7A

Manual Integrations
 APPROVED

mohammad
 8/12/2020 1:47:33 PM

Quant Time: Aug 11 16:23:50 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	6.79	152	139792	20.00	ng	-0.01
21) Naphthalene-d8	8.07	136	553626	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	284223	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	582404	20.00	ng	0.00
76) Chrysene-d12	13.96	240	534979	20.00	ng	0.00
86) Perylene-d12	15.42	264	532481	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	663623	77.82	ng	0.00
7) Phenol-d6	6.43	99	811784	73.70	ng	-0.02
23) Nitrobenzene-d5	7.36	82	520859	54.44	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	224281	72.09	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1025591	53.88	ng	-0.01
79) Terphenyl-d14	12.91	244	1126081	45.76	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	63739	3.108	ng	99
71) Phenanthrene	11.35	178	92886	3.101	ng	97
75) Fluoranthene	12.53	202	211611	6.374	ng	99
78) Pyrene	12.76	202	204007	5.394	ng	99
81) Benzo(a)anthracene	13.95	228	116906	3.609	ng	98
83) Chrysene	13.99	228	98567	2.895	ng	97
88) Benzo(b)fluoranthene	14.99	252	124554m	3.870	ng	
90) Benzo(a)pyrene	15.35	252	94405	3.215	ng	97

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

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