

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
 Data File : BF120200.D
 Acq On : 24 Apr 2020 14:18
 Operator : MA/SJ
 Sample : L2117-06 5X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-042220

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:04:24 PM

Quant Time: Apr 24 14:56:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 10:24:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	132233	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	524399	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	244336	20.00	ng	0.00
64) Phenanthrene-d10	11.18	188	393366	20.00	ng	0.00
76) Chrysene-d12	13.82	240	251544	20.00	ng	-0.01
87) Perylene-d12	15.22	264	238176	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	168585	22.03	ng	0.00
7) Phenol-d6	6.31	99	227668	23.09	ng	0.00
23) Nitrobenzene-d5	7.22	82	130214	12.85	ng	-0.02
42) 2,4,6-Tribromophenol	10.49	330	50019	16.72	ng	-0.01
45) 2-Fluorobiphenyl	9.03	172	300318	17.45	ng	0.00
79) Terphenyl-d14	12.77	244	232283	15.54	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.42	163	39400	2.085	ng	99
71) Phenanthrene	11.20	178	63127	3.015	ng	96
75) Fluoranthene	12.39	202	78395	3.471	ng	97
78) Pyrene	12.62	202	83267	3.927	ng	96
81) Benzo(a)anthracene	13.80	228	34859	2.107	ng	# 93
83) Chrysene	13.84	228	35493	2.111	ng	# 94
88) Benzo(b)fluoranthene	14.82	252	40476m	2.362	ng	

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

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