

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
 Data File : BF119622.D
 Acq On : 28 Mar 2020 17:43
 Operator : CG/JU
 Sample : L2029-05
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1F-(0-2)

Manual Integrations
 APPROVED

mohammad
 4/2/2020 9:15:04 AM

Quant Time: Mar 30 06:05:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 27 07:52:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	302002	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1123417	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	519188	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	806856	20.00	ng	0.00
76) Chrysene-d12	14.00	240	602143	20.00	ng	-0.01
87) Perylene-d12	15.47	264	524656	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1797519	94.75	ng	0.01
7) Phenol-d6	6.46	99	2353960	97.13	ng	0.00
23) Nitrobenzene-d5	7.40	82	1414528	68.43	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	410479	76.64	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	2551121	72.79	ng	0.00
79) Terphenyl-d14	12.95	244	1892668	61.58	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.59	163	212315	5.757	ng	100
71) Phenanthrene	11.39	178	230799	5.517	ng	97
75) Fluoranthene	12.57	202	215279	4.755	ng	98
78) Pyrene	12.80	202	188718	3.819	ng	95
81) Benzo(a)anthracene	13.99	228	84945	2.169	ng	95
88) Benzo(b)fluoranthene	15.04	252	83027m	2.369	ng	

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

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