

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030420\
 Data File : BF119194.D
 Acq On : 4 Mar 2020 14:55
 Operator : CG/JU
 Sample : L1762-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-06-030220

Manual Integrations
 APPROVED

mohammad
 3/5/2020 3:54:21 PM

Quant Time: Mar 04 16:34:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	138497	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	505202	20.00	ng	-0.01
39) Acenaphthene-d10	9.77	164	250427	20.00	ng	-0.01
64) Phenanthrene-d10	11.26	188	376417	20.00	ng	0.00
76) Chrysene-d12	13.90	240	267434	20.00	ng	0.00
87) Perylene-d12	15.33	264	314711	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	758788	91.56	ng	0.00
7) Phenol-d6	6.39	99	911577	90.44	ng	-0.01
23) Nitrobenzene-d5	7.30	82	595074	62.19	ng	-0.02
42) 2,4,6-Tribromophenol	10.56	330	248649	75.90	ng	-0.01
45) 2-Fluorobiphenyl	9.10	172	1132102	66.60	ng	-0.01
79) Terphenyl-d14	12.84	244	1000964	64.44	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.49	163	185905	9.708	ng	99
71) Phenanthrene	11.28	178	110354	5.376	ng	98
75) Fluoranthene	12.47	202	157005	7.205	ng	99
78) Pyrene	12.70	202	134403	6.259	ng	97
81) Benzo(a)anthracene	13.89	228	57634	3.163	ng	98
83) Chrysene	13.92	228	45381	2.499	ng	96
88) Benzo(b)fluoranthene	14.92	252	63427m	3.058	ng	
90) Benzo(a)pyrene	15.27	252	42543	2.272	ng	# 93

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

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