

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100620\
 Data File : BF121859.D
 Acq On : 6 Oct 2020 18:32
 Operator : JU/CG
 Sample : L4231-02 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 337

Manual Integrations
 APPROVED

mohammad
 10/7/2020 10:41:05 AM

Quant Time: Oct 07 09:07:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	124140	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	470726	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	210397	20.00	ng	-0.01
64) Phenanthrene-d10	11.29	188	306327	20.00	ng	0.00
76) Chrysene-d12	13.93	240	227329	20.00	ng	0.00
86) Perylene-d12	15.39	264	217799	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	130348	15.60	ng	0.00
7) Phenol-d6	6.40	99	171762	15.67	ng	-0.01
23) Nitrobenzene-d5	7.32	82	108178	12.34	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	30405	11.63	ng	-0.01
45) 2-Fluorobiphenyl	9.12	172	199946	14.39	ng	-0.01
79) Terphenyl-d14	12.86	244	140816	12.55	ng	-0.02
Target Compounds						
75) Fluoranthene	12.50	202	43439	2.438	ng	97
78) Pyrene	12.73	202	51445	3.097	ng	98
83) Chrysene	13.95	228	41860	2.874	ng	# 96
84) Bis(2-ethylhexyl)phthalate	13.91	149	62755	6.993	ng	# 87
88) Benzo(b)fluoranthene	14.97	252	41804m	2.871	ng	

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

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