

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117713.D
 Acq On : 7 Nov 2019 19:07
 Operator : JU
 Sample : K5723-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 8-C

Manual Integrations
 APPROVED

Sohil

11/8/2019 3:31:38 PM

Quant Time: Nov 08 02:32:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	284085	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1106685	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	572400	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1044307	20.00	ng	0.00
76) Chrysene-d12	14.12	240	703679	20.00	ng	-0.01
87) Perylene-d12	15.63	264	791995	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1316420	83.88	ng	0.00
7) Phenol-d6	6.55	99	1658928	85.73	ng	-0.01
23) Nitrobenzene-d5	7.49	82	963978	51.43	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	432572	77.25	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	1754335	60.47	ng	0.00
79) Terphenyl-d14	13.06	244	1799741	52.65	ng	0.00
Target Compounds						
10) Phenol	6.56	94	73608	3.530	ng	98
50) Dimethylphthalate	9.69	163	246996	6.383	ng	98
75) Fluoranthene	12.69	202	187486	3.668	ng	94
78) Pyrene	12.92	202	192829	3.754	ng	95
83) Chrysene	14.14	228	101944	2.357	ng	97
88) Benzo(b)fluoranthene	15.17	252	130731m	2.621	ng	

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

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