

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117709.D
 Acq On : 7 Nov 2019 17:19
 Operator : JU
 Sample : K5723-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 7-B

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:35 PM

Quant Time: Nov 08 02:24:46 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	324539	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1245393	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	642878	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1175127	20.00	ng	0.00
76) Chrysene-d12	14.12	240	744336	20.00	ng	-0.01
87) Perylene-d12	15.63	264	851336	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1297123	72.35	ng	0.00
7) Phenol-d6	6.55	99	1651413	74.70	ng	-0.01
23) Nitrobenzene-d5	7.49	82	978463	46.38	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	448978	71.39	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	1931177	59.27	ng	0.00
79) Terphenyl-d14	13.06	244	2032654	56.22	ng	0.00
Target Compounds						
10) Phenol	6.56	94	79767	3.349	ng	99
50) Dimethylphthalate	9.69	163	331580	7.629	ng	99
75) Fluoranthene	12.69	202	166055	2.887	ng	93
78) Pyrene	12.92	202	156292	2.876	ng	97
88) Benzo(b)fluoranthene	15.17	252	111848m	2.086	ng	

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

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