

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117710.D
 Acq On : 7 Nov 2019 17:46
 Operator : JU
 Sample : K5723-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 7-C

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:36 PM

Quant Time: Nov 08 02:26:39 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	320227	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1251519	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	657947	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1181228	20.00	ng	0.00
76) Chrysene-d12	14.12	240	720030	20.00	ng	-0.01
87) Perylene-d12	15.63	264	839820	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1736682	98.17	ng	0.00
7) Phenol-d6	6.55	99	2155061	98.80	ng	0.00
23) Nitrobenzene-d5	7.49	82	1318861	62.22	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	584219	90.77	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2342432	70.24	ng	0.00
79) Terphenyl-d14	13.06	244	2387089	68.25	ng	0.00
Target Compounds						
10) Phenol	6.56	94	103863	4.419	ng	81
50) Dimethylphthalate	9.69	163	387922	8.721	ng	98
75) Fluoranthene	12.69	202	160027	2.768	ng	94
78) Pyrene	12.92	202	155465	2.958	ng	97
88) Benzo(b)fluoranthene	15.17	252	107603m	2.034	ng	

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

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