

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117714.D
 Acq On : 7 Nov 2019 19:33
 Operator : JU
 Sample : K5723-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 8-D

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:39 PM

Quant Time: Nov 08 02:34:26 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	570265	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	2154362	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	1151250	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	2099886	20.00	ng	0.00
76) Chrysene-d12	14.12	240	1427526	20.00	ng	0.00
87) Perylene-d12	15.63	264	1662991	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1674309	53.15	ng	0.00
7) Phenol-d6	6.55	99	2119789	54.57	ng	0.00
23) Nitrobenzene-d5	7.49	82	1262890	34.61	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	603428	53.58	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2257765	38.69	ng	0.00
79) Terphenyl-d14	13.06	244	2465596	35.56	ng	0.00
Target Compounds						
10) Phenol	6.56	94	101872	2.434	ng	84
50) Dimethylphthalate	9.69	163	359627	4.621	ng	98
75) Fluoranthene	12.69	202	355535	3.459	ng	96
78) Pyrene	12.92	202	381254	3.659	ng	97
83) Chrysene	14.14	228	202869	2.312	ng	97
88) Benzo(b)fluoranthene	15.18	252	269770m	2.576	ng	

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

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