

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
 Data File : BF117732.D
 Acq On : 8 Nov 2019 15:50
 Operator : JU
 Sample : K5722-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 6-B

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:05:59 AM

Quant Time: Nov 09 02:24:08 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	336850	20.00	ng	-0.01
21) Naphthalene-d8	8.22	136	1320466	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	726241	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	1422024	20.00	ng	0.00
76) Chrysene-d12	14.12	240	1029714	20.00	ng	-0.01
87) Perylene-d12	15.62	264	912535	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1748952	93.98	ng	0.00
7) Phenol-d6	6.55	99	2259864	98.49	ng	-0.01
23) Nitrobenzene-d5	7.49	82	1362482	60.92	ng	-0.02
42) 2,4,6-Tribromophenol	10.77	330	665338	93.65	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	2560169	69.55	ng	-0.01
79) Terphenyl-d14	13.06	244	3086664	61.71	ng	0.00
Target Compounds						
10) Phenol	6.56	94	52409	2.120	ng	98
50) Dimethylphthalate	9.69	163	216899	4.418	ng	98
71) Phenanthrene	11.49	178	285994	4.196	ng	95
75) Fluoranthene	12.68	202	341208	4.903	ng	95
78) Pyrene	12.91	202	358650	4.771	ng	97
81) Benzo(a)anthracene	14.10	228	145678	2.273	ng	98
83) Chrysene	14.14	228	184706	2.918	ng	98
88) Benzo(b)fluoranthene	15.17	252	211256m	3.676	ng	

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

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