

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112719\
 Data File : BF118052.D
 Acq On : 27 Nov 2019 20:28
 Operator : JU
 Sample : K5977-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WS-112119-2-5-COMP-B5

Manual Integrations
 APPROVED

mohammad
 11/29/2019 8:53:55 AM

Quant Time: Nov 28 01:23:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	146465	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	568755	20.00	ng	-0.01
39) Acenaphthene-d10	9.93	164	287945	20.00	ng	-0.01
64) Phenanthrene-d10	11.43	188	466877	20.00	ng	0.00
76) Chrysene-d12	14.08	240	307842	20.00	ng	0.00
87) Perylene-d12	15.58	264	316765	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	733670	88.67	ng	0.00
7) Phenol-d6	6.51	99	934246	93.61	ng	-0.01
23) Nitrobenzene-d5	7.45	82	564692	59.02	ng	-0.02
42) 2,4,6-Tribromophenol	10.73	330	256441	84.12	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1107804	69.02	ng	-0.01
79) Terphenyl-d14	13.02	244	966234	57.36	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.65	163	234619	11.619	ng	98
71) Phenanthrene	11.45	178	171425	7.303	ng	96
78) Pyrene	12.87	202	174606	7.018	ng	99
81) Benzo(a)anthracene	14.07	228	87770	4.315	ng	98
83) Chrysene	14.10	228	74600	3.784	ng	97
88) Benzo(b)fluoranthene	15.14	252	99874m	4.853	ng	
90) Benzo(a)pyrene	15.52	252	65383	3.613	ng	98
92) Benzo(g,h,i)perylene	17.55	276	32438	2.091	ng	98

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

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