

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112719\
 Data File : BF118046.D
 Acq On : 27 Nov 2019 17:42
 Operator : JU
 Sample : K5959-06 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WS-112019-0-2-COMP-B4

Manual Integrations
 APPROVED

mohammad
 11/29/2019 8:53:36 AM

Quant Time: Nov 28 00:32:11 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	6.88	152	145415	20.00	ng	-0.01
21) Naphthalene-d8	8.17	136	545393	20.00	ng	-0.01
39) Acenaphthene-d10	9.93	164	274015	20.00	ng	-0.01
64) Phenanthrene-d10	11.43	188	445363	20.00	ng	0.00
76) Chrysene-d12	14.08	240	317763	20.00	ng	0.00
87) Perylene-d12	15.58	264	352765	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	316245	38.50	ng	-0.01
7) Phenol-d6	6.50	99	415731	41.96	ng	-0.02
23) Nitrobenzene-d5	7.44	82	252968	27.57	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	98493	33.95	ng	-0.02
45) 2-Fluorobiphenyl	9.25	172	540803	35.41	ng	-0.02
79) Terphenyl-d14	13.02	244	489548	28.16	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.64	163	85921	4.472	ng	99
71) Phenanthrene	11.45	178	49226	2.198	ng	98
78) Pyrene	12.87	202	94801	3.692	ng	97
81) Benzo(a)anthracene	14.07	228	63887	3.042	ng	# 91
83) Chrysene	14.10	228	65454	3.217	ng	97
88) Benzo(b)fluoranthene	15.13	252	89162m	3.890	ng	
90) Benzo(a)pyrene	15.52	252	56792	2.818	ng	# 88

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

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