

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110429.D
 Acq On : 2 Nov 2018 23:15
 Operator : JU/SJ
 Sample : J5769-03 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 WAGARAW-RD-2

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:51 AM

Quant Time: Nov 03 01:01:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	82958	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	314686	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	124849	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	201631	20.00	ng	0.00
76) Chrysene-d12	14.13	240	151985	20.00	ng	0.00
87) Perylene-d12	15.66	264	114425	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	193156	37.92	ng	0.00
7) Phenol-d6	6.57	99	241752	37.10	ng	0.00
23) Nitrobenzene-d5	7.52	82	141667	26.70	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	40313	32.60	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	247615	31.14	ng	0.00
79) Terphenyl-d14	13.07	244	193037	26.41	ng	0.00
Target Compounds						
10) Phenol	6.59	94	14854	2.133	ng	# 77
50) Dimethylphthalate	9.70	163	35803	3.885	ng	# 98
71) Phenanthrene	11.51	178	70085	6.643	ng	98
75) Fluoranthene	12.70	202	214227	20.008	ng	99
78) Pyrene	12.93	202	192566	17.673	ng	98
81) Benzo(a)anthracene	14.12	228	92262	10.237	ng	97
83) Chrysene	14.16	228	92077	10.756	ng	99
86) Indeno(1,2,3-cd)pyrene	17.22	276	41022	5.776	ng	# 94
88) Benzo(b)fluoranthene	15.20	252	100013m	15.136	ng	
89) Benzo(k)fluoranthene	15.23	252	39078m	6.016	ng	
90) Benzo(a)pyrene	15.59	252	65394	10.755	ng	97
92) Benzo(a,h,i)perylene	17.70	276	45989	8.896	ng	# 91

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

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