

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
 Data File : BF110505.D
 Acq On : 6 Nov 2018 10:43
 Operator : JU/SJ
 Sample : J5813-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-A

Manual Integrations
 APPROVED

Sohil
 11/7/2018 2:54:18 PM

Quant Time: Nov 06 12:12:12 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	6.95	152	104592	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	417196	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	188587	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	290768	20.00	ng	0.00
76) Chrysene-d12	14.13	240	182277	20.00	ng	0.00
87) Perylene-d12	15.66	264	167129	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	686480	106.88	ng	0.01
7) Phenol-d6	6.59	99	873661	106.35	ng	0.00
23) Nitrobenzene-d5	7.52	82	555346	78.95	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	168310	90.10	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	929879	77.43	ng	0.00
79) Terphenyl-d14	13.07	244	557894	63.65	ng	0.00
Target Compounds						
10) Phenol	6.59	94	59301	6.753	ng	81
50) Dimethylphthalate	9.70	163	165803	11.910	ng	99
71) Phenanthrene	11.51	178	48293	3.174	ng	99
75) Fluoranthene	12.70	202	93518	6.057	ng	99
78) Pyrene	12.93	202	75097	5.747	ng	97
81) Benzo(a)anthracene	14.12	228	30456	2.818	ng	98
83) Chrysene	14.16	228	32710	3.186	ng	95
88) Benzo(b)fluoranthene	15.20	252	43278m	4.484	ng	
90) Benzo(a)pyrene	15.59	252	26296	2.961	ng	# 93

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

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