

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\
 Data File : BF110318.D
 Acq On : 30 Oct 2018 22:06
 Operator : JU/SJ
 Sample : J5711-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4

Manual Integrations
 APPROVED

Sohil
 10/31/2018 11:15:07 AM

Quant Time: Oct 31 06:03:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 17:29:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	98754	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	396020	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	168113	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	245803	20.00	ng	0.00
76) Chrysene-d12	14.16	240	179149	20.00	ng	0.00
87) Perylene-d12	15.70	264	120401	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	602810	99.40	ng	0.00
7) Phenol-d6	6.60	99	759727	97.94	ng	0.00
23) Nitrobenzene-d5	7.53	82	460213	68.93	ng	0.00
42) 2,4,6-Tribromophenol	10.81	330	132174	79.37	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	741202	69.23	ng	0.00
79) Terphenyl-d14	13.09	244	457689	53.13	ng	0.00
Target Compounds						
10) Phenol	6.60	94	35591	4.293	ng	82
31) Naphthalene	8.27	128	53231	2.916	ng	97
37) 2-Methylnaphthalene	8.97	142	25313	2.096	ng	95
50) Dimethylphthalate	9.72	163	142215	11.460	ng	# 98
71) Phenanthrene	11.53	178	30025	2.334	ng	97
75) Fluoranthene	12.73	202	49227	3.771	ng	97
78) Pyrene	12.96	202	77164	6.008	ng	97
81) Benzo(a)anthracene	14.15	228	37124	3.494	ng	# 85
83) Chrysene	14.19	228	34877	3.456	ng	# 91
88) Benzo(b)fluoranthene	15.24	252	37664m	5.417	ng	
90) Benzo(a)pyrene	15.64	252	30358	4.745	ng	# 73
92) Benzo(a,h,i)perylene	17.76	276	15178	2.790	ng	# 76

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

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