

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
 Data File : BF107349.D
 Acq On : 12 Jul 2018 19:47
 Operator : JU/SJ
 Sample : J3924-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-1-4

Manual Integrations
 APPROVED

Sohil
 7/13/2018 11:17:15 AM

Quant Time: Jul 13 03:43:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 11 16:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	62995	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	226348	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	100921	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	181828	20.00	ng	0.00
75) Chrysene-d12	14.14	240	175131	20.00	ng	0.00
86) Perylene-d12	15.67	264	126892	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	466993	125.53	ng	0.01
7) Phenol-d6	6.58	99	569085	122.49	ng	0.00
23) Nitrobenzene-d5	7.50	82	358253	87.96	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	146214	125.00	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	591626	102.21	ng	0.00
78) Terphenyl-d14	13.06	244	604679	83.64	ng	0.00
Target Compounds						
10) Phenol	6.60	94	33376	6.295	ng	98
49) Dimethylphthalate	9.68	163	146320	19.331	ng	98
70) Phenanthrene	11.50	178	34627	3.948	ng	98
74) Fluoranthene	12.69	202	94458	9.772	ng	98
77) Pyrene	12.93	202	93584	8.103	ng	98
80) Benzo(a)anthracene	14.12	228	58603	5.490	ng	99
82) Chrysene	14.16	228	49702	5.061	ng	99
85) Indeno(1,2,3-cd)pyrene	17.26	276	25196	3.024	ng	# 88
87) Benzo(b)fluoranthene	15.21	252	58843m	7.465	ng	
88) Benzo(k)fluoranthene	15.24	252	21291m	2.836	ng	
89) Benzo(a)pyrene	15.61	252	40393	5.764	ng	97
91) Benzo(a,h,i)perylene	17.75	276	23610	4.068	ng	# 89

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

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