

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
 Data File : BF107348.D
 Acq On : 12 Jul 2018 19:20
 Operator : JU/SJ
 Sample : J3918-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-204-255-ETS

Manual Integrations
 APPROVED

Sohil
 7/13/2018 11:17:14 AM

Quant Time: Jul 13 03:39:55 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	68979	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	237510	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	100825	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	184029	20.00	ng	0.00
75) Chrysene-d12	14.14	240	192609	20.00	ng	0.00
86) Perylene-d12	15.67	264	136410	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	512211	125.74	ng	0.01
7) Phenol-d6	6.58	99	630108	123.86	ng	0.00
23) Nitrobenzene-d5	7.50	82	393050	91.97	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	154722	132.40	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	636901	111.52	ng	0.00
78) Terphenyl-d14	13.07	244	710439	89.35	ng	0.00
Target Compounds						
10) Phenol	6.60	94	37307	6.426	ng	99
49) Dimethylphthalate	9.68	163	133452	17.648	ng	98
70) Phenanthrene	11.50	178	27439	3.091	ng	99
74) Fluoranthene	12.70	202	64534	6.596	ng	93
77) Pyrene	12.93	202	65483	5.156	ng	98
80) Benzo(a)anthracene	14.12	228	39028	3.325	ng	98
82) Chrysene	14.16	228	34295	3.176	ng	97
83) Bis(2-ethylhexyl)phthalate	14.08	149	34430	5.157	ng	# 94
87) Benzo(b)fluoranthene	15.21	252	40051m	4.727	ng	
89) Benzo(a)pyrene	15.61	252	26707	3.545	ng	96
91) Benzo(g,h,i)perylene	17.75	276	16113	2.582	ng	# 89

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

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