

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
 Data File : BF107351.D
 Acq On : 12 Jul 2018 20:40
 Operator : JU/SJ
 Sample : J3927-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-RBR-200031

Manual Integrations
 APPROVED

Sohil
 7/13/2018 11:17:16 AM

Quant Time: Jul 13 03:53:22 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	65927	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	235453	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	97795	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	179987	20.00	ng	0.00
75) Chrysene-d12	14.14	240	192190	20.00	ng	0.00
86) Perylene-d12	15.67	264	131989	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	349421	89.75	ng	0.00
7) Phenol-d6	6.58	99	424305	87.27	ng	0.00
23) Nitrobenzene-d5	7.50	82	260912	61.58	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	97395	85.93	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	444390	75.22	ng	0.00
78) Terphenyl-d14	13.06	244	506298	63.81	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	76982	10.496	ng	97
74) Fluoranthene	12.69	202	38183	3.991	ng	93
77) Pyrene	12.93	202	35141	2.773	ng	98
82) Chrysene	14.16	228	22005	2.042	ng	96
87) Benzo(b)fluoranthene	15.21	252	22519m	2.747	ng	

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

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