

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072518\
 Data File : BF107669.D
 Acq On : 25 Jul 2018 20:44
 Operator : JU/SJ
 Sample : J4031-04 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PLANTER-DOC-4

Manual Integrations
 APPROVED

Sohil
 7/26/2018 2:01:58 PM

Quant Time: Jul 26 11:03:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	202379	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	798475	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	311393	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	570895	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	431608	20.00	ng	-0.01
86) Perylene-d12	15.69	264	383676	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	505847	40.19	ng	-0.01
7) Phenol-d6	6.60	99	693722	45.90	ng	-0.02
23) Nitrobenzene-d5	7.53	82	391513	33.09	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	163996	46.31	ng	-0.02
44) 2-Fluorobiphenyl	9.32	172	765929	34.64	ng	-0.02
78) Terphenyl-d14	13.08	244	612283	36.32	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.71	163	65208	2.779	ng	98
74) Fluoranthene	12.72	202	122258	4.095	ng	95
77) Pyrene	12.95	202	107463m	3.640	ng	
82) Chrysene	14.18	228	54116	2.227	ng	99
87) Benzo(b)fluoranthene	15.23	252	56361m	2.684	ng	

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

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