

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020718\
 Data File : BF102835.D
 Acq On : 8 Feb 2018 00:50
 Operator : SJ/JU
 Sample : J1473-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3

Manual Integrations
 APPROVED

Sohil
 2/8/2018 1:41:50 PM

Quant Time: Feb 08 06:44:55 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	187959	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	757016	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	301991	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	452196	20.00	ng	0.00
75) Chrysene-d12	14.05	240	349894	20.00	ng	0.00
86) Perylene-d12	15.53	264	272291	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1381116	111.59	ng	0.02
7) Phenol-d6	6.52	99	1622099	108.85	ng	0.01
23) Nitrobenzene-d5	7.45	82	1011513	92.69	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	274925	113.11	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	1517514	83.38	ng	0.00
78) Terphenyl-d14	13.00	244	1116523	75.56	ng	0.00
Target Compounds						
10) Phenol	6.52	94	39991	2.504	ng	# 63
49) Dimethylphthalate	9.63	163	93494	3.971	ng	# 98
70) Phenanthrene	11.43	178	82235	3.265	ng	98
74) Fluoranthene	12.62	202	140439	5.543	ng	99
77) Pyrene	12.85	202	114446	4.171	ng	98
80) Benzo(a)anthracene	14.04	228	44249	2.011	ng	98
82) Chrysene	14.07	228	55968m	2.523	ng	
87) Benzo(b)fluoranthene	15.09	252	54780m	3.103	ng	

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

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