

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089472.D
 Acq On : 31 Jul 2016 15:49
 Operator : UM/SJ
 Sample : H4257-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP-2

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:17:58 PM

Quant Time: Aug 01 05:40:02 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 04:11:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	45037	20.00	ng	-0.01
21) Naphthalene-d8	7.78	136	175295	20.00	ng	-0.01
38) Acenaphthene-d10	9.52	164	86459	20.00	ng	-0.01
63) Phenanthrene-d10	10.99	188	149674	20.00	ng	-0.01
75) Chrysene-d12	13.61	240	82008	20.00	ng	-0.01
86) Perylene-d12	14.94	264	82397	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	374449	132.76	ng	0.00
7) Phenol-d6	6.16	99	556329	149.25	ng	-0.01
23) Nitrobenzene-d5	7.06	82	311464	104.88	ng	-0.02
41) 2,4,6-Tribromophenol	10.31	330	95392	133.22	ng	-0.01
44) 2-Fluorobiphenyl	8.86	172	511136	86.76	ng	-0.01
78) Terphenyl-d14	12.57	244	341480	97.46	ng	-0.01
Target Compounds						
10) Phenol	6.17	94	8945	2.17	ng	Qvalue # 77
49) Dimethylphthalate	9.26	163	94791	14.41	ng	98
74) Fluoranthene	12.19	202	16710	2.06	ng	96
77) Pyrene	12.42	202	15569	2.50	ng	97
80) Benzo(a)anthracene	13.60	228	10033	2.17	ng	97
85) Indeno(1,2,3-cd)pyrene	16.13	276	10028	2.86	ng	# 91
87) Benzo(b)fluoranthene	14.61	252	14123m	2.76	ng	
89) Benzo(a)pyrene	14.89	252	13456	2.93	ng	97
91) Benzo(g,h,i)perylene	16.47	276	10011	2.82	ng	95

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

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