

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
 Data File : BF089385.D
 Acq On : 29 Jul 2016 2:06
 Operator : UM/SJ
 Sample : H4216-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3(LOT109)0-1

Quant Time: Jul 29 03:30:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 16:02:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	56117	20.00	ng	-0.02
21) Naphthalene-d8	7.82	136	212509	20.00	ng	-0.01
38) Acenaphthene-d10	9.55	164	98090	20.00	ng	-0.02
63) Phenanthrene-d10	11.03	188	148129	20.00	ng	-0.02
75) Chrysene-d12	13.65	240	113544	20.00	ng	-0.02
86) Perylene-d12	14.98	264	101805	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.11	112	318842	90.72	ng	0.00
7) Phenol-d6	6.18	99	410414	88.36	ng	-0.02
23) Nitrobenzene-d5	7.10	82	262409	72.89	ng	-0.02
41) 2,4,6-Tribromophenol	10.34	330	64579	79.49	ng	-0.02
44) 2-Fluorobiphenyl	8.89	172	447716	66.98	ng	-0.02
78) Terphenyl-d14	12.60	244	280010	57.72	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.29	163	139554	18.69	ng	98
70) Phenanthrene	11.05	178	39757	5.20	ng	99
74) Fluoranthene	12.23	202	59990	7.46	ng	98
77) Pyrene	12.46	202	59380	6.90	ng	99
80) Benzo(a)anthracene	13.63	228	25964	4.05	ng	# 82
82) Chrysene	13.67	228	22656m	3.32	ng	
85) Indeno(1,2,3-cd)pyrene	16.18	276	25720	5.30	ng	93
87) Benzo(b)fluoranthene	14.64	252	41252m	6.52	ng	
88) Benzo(k)fluoranthene	14.65	252	13078m	2.25	ng	
89) Benzo(a)pyrene	14.94	252	28939	5.10	ng	95
91) Benzo(g,h,i)perylene	16.53	276	26042	5.93	ng	97

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

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