

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079727.D
 Acq On : 5 Jun 2015 5:12
 Operator : TP/IZ
 Sample : G2474-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-6

Quant Time: Jun 06 04:01:24 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 06 03:31:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	102205	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	399346	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	208989	20.00	ng	0.00
63) Phenanthrene-d10	12.09	188	424258	20.00	ng	0.01
75) Chrysene-d12	15.02	240	465817	20.00	ng	0.03
86) Perylene-d12	16.96	264	464766	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	644318	108.10	ng	0.00
7) Phenol-d6	7.11	99	946945	122.04	ng	0.01
23) Nitrobenzene-d5	8.07	82	470555	71.74	ng	0.00
41) 2,4,6-Tribromophenol	11.37	330	224744	119.75	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	1003260	69.50	ng	0.00
78) Terphenyl-d14	13.76	244	1238658	62.65	ng	0.02
Target Compounds						
49) Dimethylphthalate	10.26	163	134236	8.34	ng	Qvalue # 98
70) Phenanthrene	12.11	178	81320	3.30	ng	97
74) Fluoranthene	13.36	202	187123	6.99	ng	95
77) Pyrene	13.62	202	185097	6.39	ng	98
80) Benzo(a)anthracene	15.01	228	103219	3.62	ng	97
82) Chrysene	15.05	228	93993	3.53	ng	98
85) Indeno(1,2,3-cd)pyrene	18.96	276	71749	2.39	ng	94
87) Benzo(b)fluoranthene	16.38	252	131995m	4.40	ng	
89) Benzo(a)pyrene	16.87	252	92640	3.38	ng	# 93
91) Benzo(g,h,i)perylene	19.57	276	70417	2.72	ng	97

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

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