

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079728.D
 Acq On : 5 Jun 2015 5:41
 Operator : TP/IZ
 Sample : G2474-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-7

Quant Time: Jun 06 04:06:08 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	63113	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	244634	20.00	ng	0.00
38) Acenaphthene-d10	10.57	164	123074	20.00	ng	-0.01
63) Phenanthrene-d10	12.09	188	263547	20.00	ng	0.01
75) Chrysene-d12	15.02	240	300864	20.00	ng	0.03
86) Perylene-d12	16.96	264	307750	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	337335	91.65	ng	0.00
7) Phenol-d6	7.10	99	408804	85.32	ng	0.00
23) Nitrobenzene-d5	8.06	82	211004	52.52	ng	-0.01
41) 2,4,6-Tribromophenol	11.37	330	113621	102.81	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	579403	68.16	ng	0.00
78) Terphenyl-d14	13.76	244	723002	56.62	ng	0.02
Target Compounds						Qvalue
49) Dimethylphthalate	10.26	163	61342	6.47	ng	98
70) Phenanthrene	12.11	178	47859	3.12	ng	96
74) Fluoranthene	13.36	202	122815	7.39	ng	96
77) Pyrene	13.62	202	115216	6.16	ng	97
80) Benzo(a)anthracene	15.01	228	63094	3.43	ng	97
82) Chrysene	15.05	228	59731	3.47	ng	99
85) Indeno(1,2,3-cd)pyrene	18.95	276	48393	2.49	ng	# 93
87) Benzo(b)fluoranthene	16.38	252	88452m	4.46	ng	
89) Benzo(a)pyrene	16.87	252	60405	3.33	ng	# 95
91) Benzo(g,h,i)perylene	19.57	276	50736	2.96	ng	96

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

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