

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079726.D
 Acq On : 5 Jun 2015 4:43
 Operator : TP/IZ
 Sample : G2443-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW-027

Quant Time: Jun 06 03:56:47 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	84110	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	336236	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	168585	20.00	ng	0.00
63) Phenanthrene-d10	12.09	188	358929	20.00	ng	0.01
75) Chrysene-d12	15.06	240	374949	20.00	ng	0.08
86) Perylene-d12	17.03	264	381528	20.00	ng	0.11
System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	584937	119.25	ng	0.00
7) Phenol-d6	7.11	99	834616	130.70	ng	0.01
23) Nitrobenzene-d5	8.07	82	419681	76.00	ng	0.00
41) 2,4,6-Tribromophenol	11.37	330	203731	134.57	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	933420	80.16	ng	0.00
78) Terphenyl-d14	13.78	244	1120876	70.43	ng	0.05
Target Compounds						Qvalue
48) Acenaphthylene	10.43	152	44328	2.35	ng	98
49) Dimethylphthalate	10.26	163	274717	21.15	ng	98
70) Phenanthrene	12.11	178	344531	16.50	ng	99
71) Anthracene	12.17	178	98840	4.78	ng	99
74) Fluoranthene	13.38	202	598348	26.42	ng	98
77) Pyrene	13.65	202	581065	24.94	ng	98
80) Benzo(a)anthracene	15.05	228	324019	14.14	ng	97
82) Chrysene	15.10	228	280407	13.07	ng	97
83) Bis(2-ethylhexyl)phthalate	15.01	149	29976	2.13	ng	98
85) Indeno(1,2,3-cd)pyrene	19.04	276	196689	8.13	ng	95
87) Benzo(b)fluoranthene	16.45	252	385986m	15.68	ng	
88) Benzo(k)fluoranthene	16.48	252	135449m	5.77	ng	
89) Benzo(a)pyrene	16.95	252	291725	12.97	ng	99
90) Dibenzo(a,h)anthracene	19.05	278	51949	2.50	ng	# 91
91) Benzo(g,h,i)perylene	19.65	276	196432	9.25	ng	99

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

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