

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
 Data File : BF087095.D
 Acq On : 16 May 2016 21:48
 Operator : UM/SJ
 Sample : H3057-01DL 10X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E-10(8.5-9)DL

Manual Integrations
 APPROVED

umangi
 5/17/2016 7:16:24 PM

Quant Time: May 17 13:37:44 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 17 01:12:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	166081	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	690339	20.00	ng	0.00
38) Acenaphthene-d10	9.64	164	282953	20.00	ng	0.00
63) Phenanthrene-d10	11.12	188	454395	20.00	ng	0.00
75) Chrysene-d12	13.75	240	378979	20.00	ng	-0.01
86) Perylene-d12	15.11	264	318969	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.20	112	91404	9.32	ng	0.00
7) Phenol-d6	6.27	99	129443	9.88	ng	-0.01
23) Nitrobenzene-d5	7.18	82	80340	5.84	ng	-0.01
41) 2,4,6-Tribromophenol	10.43	330	21244	7.09	ng	-0.01
44) 2-Fluorobiphenyl	8.97	172	126896	7.54	ng	-0.01
78) Terphenyl-d14	12.70	244	80757	4.52	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	7.92	128	124037	3.59	ng	100
37) 2-Methylnaphthalene	8.60	142	51889	2.57	ng	94
51) Acenaphthene	9.68	154	101611	6.30	ng	98
54) Dibenzofuran	9.85	168	125667	6.07	ng	98
57) Fluorene	10.19	166	137089	7.76	ng	97
70) Phenanthrene	11.15	178	1225033	50.52	ng	99
71) Anthracene	11.20	178	410182	16.54	ng	99
72) Carbazole	11.36	167	81744	3.83	ng	98
74) Fluoranthene	12.33	202	1111215	44.55	ng	97
77) Pyrene	12.56	202	1077865	37.15	ng	100
80) Benzo(a)anthracene	13.74	228	481125	22.61	ng	99
82) Chrysene	13.77	228	336504m	15.55	ng	
85) Indeno(1,2,3-cd)pyrene	16.35	276	145553	8.08	ng	93
87) Benzo(b)fluoranthene	14.74	252	363059m	17.51	ng	
88) Benzo(k)fluoranthene	14.75	252	149055m	8.78	ng	
89) Benzo(a)pyrene	15.06	252	289724	16.20	ng	97
90) Dibenzo(a,h)anthracene	16.37	278	32338m	2.15	ng	
91) Benzo(g,h,i)perylene	16.73	276	130571	8.53	ng	96

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

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