

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070617\
 Data File : BF096529.D
 Acq On : 6 Jul 2017 13:26
 Operator : SJ/MA
 Sample : I4053-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-070317-B

Manual Integrations
 APPROVED

mohammad
 7/7/2017 1:29:55 PM

Quant Time: Jul 06 18:56:31 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 05 13:27:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	127289	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	511549	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	213155	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	311083	20.00	ng	0.00
75) Chrysene-d12	13.87	240	225046	20.00	ng	0.00
86) Perylene-d12	15.28	264	194275	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	778651	90.29	ng	0.00
7) Phenol-d6	6.35	99	944280	89.06	ng	0.00
23) Nitrobenzene-d5	7.27	82	537633	63.15	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	144345	86.69	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	947989	76.04	ng	0.00
78) Terphenyl-d14	12.82	244	558664	53.05	ng	0.00

Target Compounds					Qvalue
10) Phenol	6.36	94	30394	2.516 ng	93
31) Naphthalene	8.02	128	56842	2.354 ng	99
37) 2-Methylnaphthalene	8.71	142	59978	3.902 ng	94
48) Acenaphthylene	9.60	152	92892	4.307 ng	97
49) Dimethylphthalate	9.47	163	240303	15.102 ng	98
51) Acenaphthene	9.78	154	48086	3.617 ng	99
54) Dibenzofuran	9.95	168	41659	2.374 ng	# 90
57) Fluorene	10.29	166	113748	9.176 ng	99
70) Phenanthrene	11.26	178	389039	23.134 ng	99
71) Anthracene	11.30	178	121231	7.074 ng	97
74) Fluoranthene	12.44	202	375624	22.782 ng	99
77) Pyrene	12.67	202	446033	24.692 ng	99
80) Benzo(a)anthracene	13.86	228	212748m	16.325 ng	
82) Chrysene	13.89	228	214709	15.733 ng	97
85) Indeno(1,2,3-cd)pyrene	16.65	276	177148	16.445 ng	95
87) Benzo(b)fluoranthene	14.88	252	330549m	27.154 ng	
88) Benzo(k)fluoranthene	14.90	252	91539m	8.406 ng	
89) Benzo(a)pyrene	15.22	252	265918	24.468 ng	96
90) Dibenz(a,h)anthracene	16.66	278	43049	5.035 ng	# 80
91) Benzo(g,h,i)perylene	17.06	276	162540	18.346 ng	97

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

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